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ITERATIVE AND DIRECT METHODS FOR SOLVING  
POISSON'S EQUATION AND THEIR  
ADAPTABILITY TO ILLIAC IV

by

James H. Ericksen

December 20, 1972





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## ABSTRACT

This paper examines iterative and direct methods for solving Poisson's equation with regard to their adaptation to ILLIAC IV. SOR, SLOR, ADI, and FACR are programmed in GLYPNIR. Detailed suggestions on ASK code for these methods are also supplied.

FACR, Fourier Analysis and Cyclic Reduction, is the fastest method on rectangular meshes. SOR, Successive Over Relaxation, seems to be the most promising for nonrectangular meshes. The methods are between thirty and forty-five times faster on ILLIAC IV than on a serial machine with speed equal to one of the ILLIAC IV PEs (Processing Elements).



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# 1. POISSON'S EQUATION IN MATRIX FORM

We will consider Poisson's equation with Dirichlet boundary conditions defined on a rectangular region. This equation can be written as

$$\psi(X,Y) = \frac{\partial^2 U(X,Y)}{\partial Y^2} + \frac{\partial^2 U(X,Y)}{\partial X^2} \quad (1)$$

where the value of  $U$  is given at the boundaries  $X = 0$ ,  $Y = 0$ ,  $X = (m-1)h_x$  and  $Y = (n-1)h_y$ . We denote  $U((i-1)h_x, (j-1)h_y)$  by  $U_{i,j}$ ,  $\psi((i-1)h_x, (j-1)h_y)$  by  $\psi_{i,j}$  and error of order  $p$  by  $O(p)$ . By dividing the region into mesh points such that the distance between mesh points in the vertical direction is  $h_y$  and in the horizontal direction is  $h_x$  we get Figure 1.a. The region in Figure 1.a can be rotated  $90^\circ$  clockwise to obtain the mesh  $U$  in Figure 1.b where  $U_{i,j}$  is the mesh point in the  $i^{\text{th}}$  row and the  $j^{\text{th}}$  column.

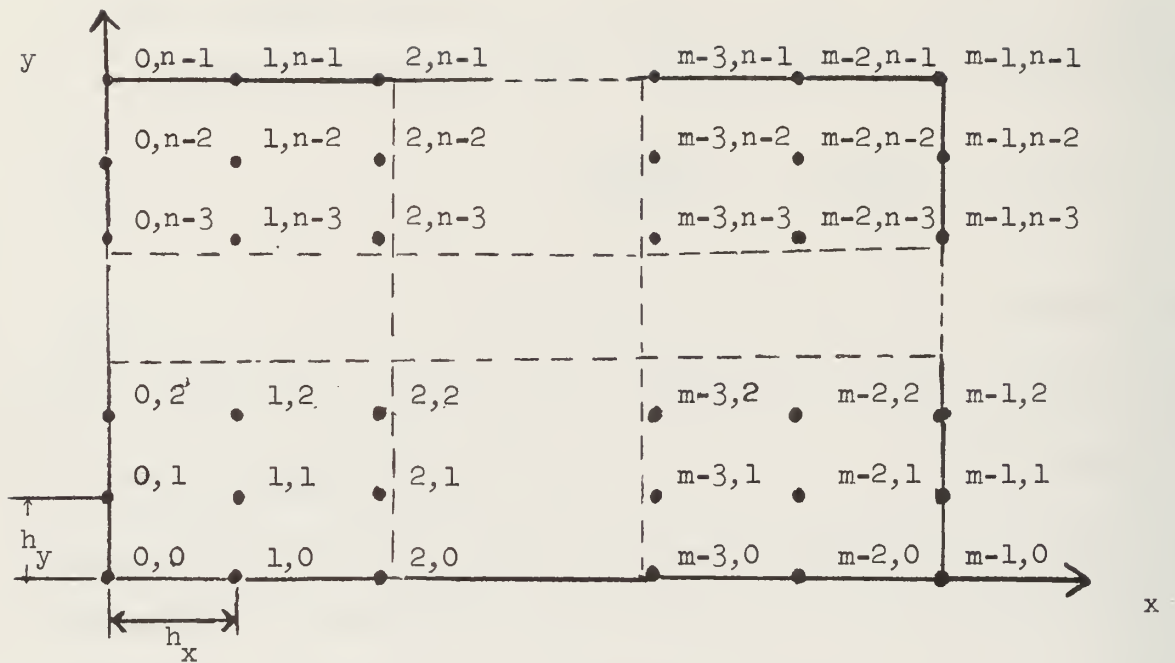
By use of Taylor's Theorem we obtain:

$$U_{i+1,j} = U_{i,j} + h_x \frac{\partial U}{\partial X}(i,j) + \frac{h_x^2}{2} \frac{\partial^2 U}{\partial X^2}(i,j) + \frac{h_x^3}{3!} \frac{\partial^3 U}{\partial X^3}(i,j) + \frac{h_x^4}{4!} \frac{\partial^4 U}{\partial X^4}(i,j) + \dots$$

$$U_{i-1,j} = U_{i,j} - h_x \frac{\partial U}{\partial X}(i,j) + \frac{h_x^2}{2} \frac{\partial^2 U}{\partial X^2}(i,j) - \frac{h_x^3}{3!} \frac{\partial^3 U}{\partial X^3}(i,j) + \frac{h_x^4}{4!} \frac{\partial^4 U}{\partial X^4}(i,j) + \dots$$

$$U_{i,j+1} = U_{i,j} + h_y \frac{\partial U}{\partial Y}(i,j) + \frac{h_y^2}{2} \frac{\partial^2 U}{\partial Y^2}(i,j) + \frac{h_y^3}{3!} \frac{\partial^3 U}{\partial Y^3}(i,j) + \frac{h_y^4}{4!} \frac{\partial^4 U}{\partial Y^4}(i,j) + \dots$$

$$U_{i,j-1} = U_{i,j} - h_y \frac{\partial U}{\partial Y}(i,j) + \frac{h_y^2}{2} \frac{\partial^2 U}{\partial Y^2}(i,j) - \frac{h_y^3}{3!} \frac{\partial^3 U}{\partial Y^3}(i,j) + \frac{h_y^4}{4!} \frac{\partial^4 U}{\partial Y^4}(i,j) + \dots$$



a) The rectangular region with the mesh points.

$U_{1,1'}$	$U_{1,2'}$	$U_{1,3'}$	$\dots$	$U_{1,n-1'}$	$U_{1,n}$
$U_{2,1'}$	$U_{2,2'}$	$U_{2,3'}$	$\dots$	$U_{2,n-1'}$	$U_{2,n}$
$U_{3,1'}$	$U_{3,2'}$	$U_{3,3'}$	$\dots$	$U_{3,n-1'}$	$U_{3,n}$
$\vdots$	$\vdots$	$\vdots$	$\dots$	$\vdots$	$\vdots$
$U_{m-1,1'}$	$U_{m-1,2'}$	$U_{m-1,3'}$	$\dots$	$U_{m-1,n-1'}$	$U_{m-1,n}$
$U_{m,1'}$	$U_{m,2'}$	$U_{m,3'}$	$\dots$	$U_{m,n-1'}$	$U_{m,n}$

b) The mesh points with separation of boundary and interior mesh points.

Figure 1. The Mesh



$$\text{Thus } \frac{\delta^2 U}{\delta Y^2}(i,j) = \frac{1}{h_y^2} (U_{i,j+1} + U_{i,j-1} - 2U_{i,j}) + O(h_y^2) \quad (2)$$

$$\text{and } \frac{\delta^2 U}{\delta X^2}(i,j) = \frac{1}{h_x^2} (U_{i+1,j} + U_{i-1,j} - 2U_{i,j}) + O(h_x^2) \quad (3)$$

Substituting (2) and (3) into (1) we obtain

$$\frac{1}{h_y^2} (U_{i,j+1} + U_{i,j-1} - 2U_{i,j}) + \frac{1}{h_x^2} (U_{i+1,j} + U_{i-1,j} - 2U_{i,j}) \approx \psi_{i,j} \quad (4)$$

This is the equation for any interior mesh point  $U_{i,j}$ . The values of the mesh points on the boundary are given, and require no equations. Let us take all the equations for the interior mesh points and place the constant terms on the right hand side of the equations. For example, the equation for  $U_{2,j}$  where  $3 \leq j \leq n-2$  will be

$$\frac{1}{h_y^2} (U_{2,j+1} + U_{2,j-1} - 2U_{2,j}) + \frac{1}{h_x^2} (U_{3,j} - 2U_{2,j}) \approx \psi_{2,j} - \frac{1}{h_x^2} U_{1,j}$$

The equation for  $U_{2,2}$  will be

$$\frac{1}{h_y^2} (U_{2,3} - 2U_{2,2}) + \frac{1}{h_x^2} (U_{3,2} - 2U_{2,2}) \approx \psi_{2,2} - \frac{1}{h_x^2} U_{1,2} - \frac{1}{h_y^2} U_{2,1}$$

Combining the constant terms into one term,  $M_{i,j}$ , for each  $U_{i,j}$  we obtain

$$\begin{aligned} M_{i,j} = & \psi_{i,j} - \frac{1}{h_y^2} (U_{i,j+1} \delta_{j+1,n} + U_{i,j-1} \delta_{j-1,1}) - \frac{1}{h_x^2} (U_{i+1,j} \delta_{i+1,m} \\ & + U_{i-1,j} \delta_{i-1,1}) \end{aligned} \quad (5)$$

$$\text{where } \delta_{k,\ell} = \begin{cases} 0 & \text{if } k \neq \ell \\ 1 & \text{if } k = \ell \end{cases}.$$

Consider the internal mesh points as a vector  $U_I$  in which the order of the mesh points is as follows:

- 1) The first interior point in the first row becomes the first element of the vector.
- 2) All interior points of row  $i$  are placed in the vector in the same order they appear in row  $i$ .
- 3) The  $(n-1)^{th}$  element of the  $i^{th}$  row is followed by the first interior point of the following row, where  $2 \leq i \leq m-2$ .

The transpose of  $U_I$  equals

$$(U_{2,2}, U_{2,3}, \dots, U_{2,n-1}; U_{3,2}, U_{3,3}, \dots, U_{3,n-1}; \dots; U_{m-1,2}, U_{m-1,3}, \dots, U_{m-1,n-1}).$$

Using this order we assemble the interior mesh point equations and consider them as the matrix equation  $AU_I = M$  where  $M$  is the vector of constant terms and  $A$  is a matrix made up of the coefficients of the mesh points.

The matrix  $A$  is a  $(m-2)(n-2)$  by  $(m-2)(n-2)$  block tridiagonal matrix of the following form:

$$A = \begin{bmatrix} B & C & & & \\ C & B & C & & \\ & C & B & C & \\ & & & & \text{Circle} \\ \text{Circle} & & & C & B & C \\ & & & & C & B \end{bmatrix} \quad (6)$$

where  $C = aI_{n-2}$ ,  $a = \frac{+1}{h_x^2}$

and  $B$  is a  $(n-2)$  by  $(n-2)$  tridiagonal matrix of the form

$$B = \begin{bmatrix} e & d & & & \\ d & e & d & & \\ & d & e & d & \\ & & d & e & d \\ & & & d & e \end{bmatrix} \quad (6.1)$$

where  $d = \frac{1}{h_y^2}$  and  $e = -(2a + 2d)$ .

There are two general approaches used in solving this system of linear equations, direct methods and iterative methods.\* In this study we have limited our discussion to three popular iterative methods and one direct method: the successive over-relaxation method (SOR), the successive line over-relaxation method (SLOR), the alternating direction implicit method (ADI), and Hockney's direct method (FACR). In the discussions of FACR three types of boundary conditions will be examined: periodic, Neumann and Dirichlet. The discussion of iterative methods considers only Dirichlet boundary conditions.

---

\* No matter which method is used in solving the equations the lower bound on the error is  $O(\max(h_y^2, h_x^2))$ . This is due to the error terms in (2) and (3).

## 2. IMPLEMENTATION OF ITERATIVE METHODS ON ILLIAC IV

### 2.1 Introduction to SLOR and SOR

Iterative methods have one thing in common. They each begin with an initial guess to the solution and improve<sup>\*</sup> upon the guess with each iteration. Thus the following notation is useful: The value of  $U_I$  after the  $\ell^{\text{th}}$  iteration is denoted by  $U_I^{(\ell)}$ .  $U_I^{(0)}$  is the initial guess and if the method converges then  $\lim_{\ell \rightarrow \infty} (U_I^{(\ell)} - U_I^{(\ell+1)}) = 0$ .

Matrix  $A$  can be divided into three matrices such that  $A = D - E - F$  where  $D$  is block diagonal,  $E$  is strictly lower triangular and  $F$  is strictly upper triangular. Thus  $AU_I = M$  can be written as  $DU_I - (E + F)U_I = M$ .

From this we get the iterative method:

$$U_I^{(\ell+1)} = D^{-1} (E + F) U_I^{(\ell)} + D^{-1} M \quad (7)$$

The matrix equation can also be grouped to obtain

$$U_I^{(\ell+1)} = (D-E)^{-1} F U_I^{(\ell)} + (D-E)^{-1} M \quad (8)$$

The asymptotic convergence rate of (8) is twice as fast as (7) [Todd, Page 391]. This is due to the fact that the matrix has Young's property A [Wachspress, Page 102]. (8) is a successive method while (7) is a simultaneous method. \*\*

By defining  $J = \frac{1}{\omega} (D - \omega E)$  and  $H = \frac{1}{\omega} (\omega F + (1-\omega)D)$  for  $\omega \neq 0$  we get  $\omega J U_I = \omega H U_I + \omega M$  from  $AU_I = M = (J-H)U_I$ .

---

\* Improvement is measured by a reduction in the error vector.

\*\* Successive methods use new information as soon as it is available while simultaneous methods use old values for the entire iteration.

We can write  $\omega JU_I = \omega HU_I + \omega M$  as

$$(D - \omega E)U_I^{(\ell+1)} = (\omega F + (1 - \omega)D)U_I^{(\ell)} + \omega M \quad (9)$$

Note that (8) and (9) are equivalent if  $\omega = 1$ . When  $0 < \omega < 2$  and  $A$  is positive definite, (9) converges. If  $1 < \omega < 2$ , then (9) is referred to as an over-relaxation method [Forsythe and Wasow, Page 261]. With different values of  $\omega$ , the convergence rate of (9) is altered. For a single relaxation parameter, it can be shown that  $\omega_b = \frac{2}{1 + \sqrt{1 - \rho^2(\beta)}}$ , where  $\beta = D^{-1}(E+F)$ , the Jacobi matrix, and  $\rho(\beta)$  is the spectral radius of  $\beta$ , gives the fastest rate of convergence [Young, Page 169].

There are two common ways in which equations (7 - 9) are implemented. The first gives point iterative methods by letting  $D$  equal the diagonal of  $A$ . Thus (7) becomes Jacobi's method, (8) becomes the formula for successive point relaxation iterative method (SR) and (9) becomes the successive point over relaxation iterative method (SOR).

If we let

$$D = \begin{bmatrix} B & \text{---} \\ \text{---} & B \end{bmatrix}$$

where  $B$  is defined in (6.1) then we will get line iterative methods.

(7) becomes the simultaneous line iterative method, (8) becomes the

successive line relaxation iterative method (SLR) and (9) becomes the successive line over-relaxation method (SLOR) [Varga, Page 199].

To give the reader a better understanding of the way these methods work, we write the individual equations for the interior mesh points  $U_{i,j}$ , where  $3 \leq i \leq m-2$  and  $3 \leq j \leq n-2$  for each method.\* First we define  $h$ ,  $v$ , and  $g$  as follows:

$$h = \frac{h_y^2}{2(h_x^2 + h_y^2)}, \quad v = \frac{h_x^2}{2(h_x^2 + h_y^2)}, \quad \text{and} \quad g = \frac{h_x^2}{2(h_x^2 + h_y^2)} \cdot \frac{h_y^2}{2(h_x^2 + h_y^2)}.$$

#### Jacobi's method (simultaneous point iterative method)

$$U_{i,j}^{(\ell+1)} = v (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell)}) + h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)}) - g M_{i,j}. \quad (10)$$

#### Simultaneous line iterative method

$$U_{i,j}^{(\ell+1)} = v (U_{i,j+1}^{(\ell+1)} + U_{i,j-1}^{(\ell+1)}) + h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)}) - g M_{i,j}.$$

#### Successive point relaxation method (SR)

$$U_{i,j}^{(\ell+1)} = v (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell+1)}) + h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell+1)}) - g M_{i,j}.$$

---

\* Note we have excluded the first row, last row, first column, and the last column of the interior mesh points to avoid boundary conditions. These points will be discussed in another section.

### Successive line relaxation method (SLR)

$$U_{i,j}^{(\ell+1)} = v (U_{i,j+1}^{(\ell+1)} + U_{i,j-1}^{(\ell+1)}) + h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)}) - g M_{i,j} . \quad (11)$$

### Successive point over-relaxation (SOR)

$$U_{i,j}^{(\ell+1)} = \omega v (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell)}) + \omega h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)}) - \omega g M_{i,j} \\ + (1-\omega) U_{i,j}^{(\ell)} . \quad (12)$$

### Successive line over-relaxation (SLOR)

$$U_{i,j}^{(\ell+1)} = v (U_{i,j+1}^{(\ell+1)} + U_{i,j-1}^{(\ell+1)}) + \omega h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)}) - \omega g M_{i,j} \\ + (1-\omega) (U_{i,j}^{(\ell)} - v (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell)})) . \quad (13)$$

## 2.2 Introduction to ADI

Let us look at the equation we are solving once again,

$$\frac{\partial^2 U(X,Y)}{\partial X^2} + \frac{\partial^2 U(X,Y)}{\partial Y^2} = \Psi(X,Y) \quad (1)$$

as before we use Taylor's Theorem and we get

$$\frac{\partial^2 U}{\partial Y^2}(i,j) = \frac{1}{h_y^2} (U_{i,j-1} + U_{i,j+1} - 2U_{i,j}) + O(h_y^2) \quad (2)$$

and

$$\frac{\partial^2 U}{\partial X^2}(i,j) = \frac{1}{h_x^2} (U_{i-1,j} + U_{i+1,j} - 2U_{i,j}) + O(h_x^2) \quad (3)$$



Thus when we form the matrix equation  $AU_I = M$  where  $A$  is the matrix (6), we are considering a horizontal part of the equation and a vertical part of the equation. We could divide  $A$  into two matrices  $H$  and  $V$  such that  $H + V = A$  where  $V$  is the matrix containing the coefficients of the vertical equations (2). Both  $H$  and  $V$  can be placed in tridiagonal form with the appropriate permutation matrix.

We can take the equation

$$AU_I = HU_I + VU_I = M \text{ and obtain}$$

$$(H - \omega_{\ell H} I)U_I = M - (V + \omega_{\ell H} I)U_I \text{ and}$$

$$(V - \omega_{\ell V} I)U_I = M - (H + \omega_{\ell V} I)U_I .$$

Thus we can get the Peaceman-Rachford ADI scheme [Varga, Page 212]. \*

$$U_I^{(\ell+1/2)} = (H - \omega_{\ell H} I)^{-1} [M - (V + \omega_{\ell H} I)U_I^{(\ell)}] \quad (14.1)$$

$$U_I^{(\ell+1)} = (V - \omega_{\ell V} I)^{-1} [M - (H + \omega_{\ell V} I)U_I^{(\ell+1/2)}] \quad (14.2)$$

The parameters,  $\omega_{\ell H}$  and  $\omega_{\ell V}$  are defined in [Wachspress, Page 192].

It will help to have the corresponding equations for the individual mesh points,  $U_{i,j}$ , where  $3 \leq i \leq n-2$  and  $3 \leq j \leq m-2$ .

$$\begin{aligned} U_{i,j}^{(\ell+1/2)} = & \frac{-1}{\omega_{\ell H} + \frac{2}{h_x^2}} \left[ M_{i,j} - \frac{1}{h_y^2} (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell)}) - \left( \omega_{\ell H} - \frac{2}{h_y^2} \right) U_{i,j}^{(\ell)} \right. \\ & \left. - \frac{1}{h_x^2} (U_{i-1,j}^{(\ell+1/2)} + U_{i+1,j}^{(\ell+1/2)}) \right]. \end{aligned} \quad (15.1)$$

---

\* ADI requires that  $A$  be positive definite. This can be obtained by using  $-AU_I = -M$  or by using  $-\omega_{\ell V}$  and  $-\omega_{\ell H}$  instead of  $\omega_{\ell V}$  and  $\omega_{\ell H}$ . We use the latter technique.

$$U_{i,j}^{(\ell+1)} = \frac{-1}{(\omega_{\ell V} + \frac{2}{h_y^2})} [M_{i,j} - \frac{1}{h_x^2} (U_{i+1,j}^{(\ell+1/2)} + U_{i-1,j}^{(\ell+1/2)}) - (\omega_{\ell V} - \frac{2}{h_x^2}) U_{i,j}^{(\ell+1/2)} - \frac{1}{h_y^2} (U_{i,j-1}^{(\ell+1)} + U_{i,j+1}^{(\ell+1)})]. \quad (15.2)$$

Examining (15.1) and (15.2) we note a strong resemblance to simultaneous column relaxation followed by simultaneous row relaxation.

### 2.3 Implementation of SOR on ILLIAC IV

#### 2.3.1 A Parallel Processor's Effect on the Algorithm

A parallel machine allows operations to be performed simultaneously which are done in separate time intervals on a conventional machine. This is referred to as overlap. The amount of overlap is dependent on the way in which the algorithm is programmed. The SOR algorithm supplies an illustration of how the flow of an algorithm is changed when programmed to maximize overlap on a parallel machine.

Let us examine the equation for SOR closer:

$$U_{i,j}^{(\ell+1)} = \omega_v (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell+1)}) + \omega_h (U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell+1)}) - \omega_g M_{i,j} + (1-\omega) U_{i,j}^{(\ell)} \quad (12)$$

for  $3 \leq i \leq m-2$  and  $3 \leq j \leq n-2$ . Note the interior points which are adjacent to one or more boundary points are not included. Also note that for the above mesh points  $M_{i,j} = \Psi_{i,j}$ . The equation for  $U_{2,2}$  would be

$$U_{2,2}^{(\ell+1)} = \omega_v U_{2,3}^{(\ell)} + \omega_h U_{3,2}^{(\ell)} - \omega_g M_{2,2} + (1-\omega) U_{2,2}^{(\ell)} \quad (16)$$

where

$$M_{2,2} = \Psi_{2,2} - \frac{1}{h_x^2} U_{1,2} - \frac{1}{h_y^2} U_{2,1}.$$

Thus

$$g M_{2,2} = g \Psi_{2,2} - v U_{2,1} - h U_{1,2}.$$

This enables us to rewrite (16) as

$$U_{2,2}^{(\ell+1)} = \omega_v (U_{2,3}^{(\ell)} + U_{2,1}^{(\ell+1)}) + \omega_h (U_{3,2}^{(\ell)} + U_{1,2}^{(\ell+1)}) - \omega_g \Psi_{2,2} \quad (17) *$$

To calculate  $U_{2,2}$  using (17) takes two more additions than using (16) but  $M_{2,2}$  is needed to use (16). On a serial machine, one would use equation (16) to improve  $U_{2,2}$ . The choice is not so straightforward on a parallel machine if when  $U_{2,2}$  is calculated in one processing element (PE),  $U_{i,j}$  is calculated in another PE where  $3 \leq i \leq m-2$  and (or)  $3 \leq j \leq n-2$ . In this case to calculate  $U_{2,2}$  using (16) we would have to turn the appropriate PE off for two additions. By using (17), we would not have to turn off PEs for those additions and would also save the calculation of  $M_{2,2}$ . A similar argument can be given for all points adjacent to boundary points. Thus the choice of the equation to evaluate a point adjacent to the boundary in one PE depends on what is being done in the other PEs.

### 2.3.2 Parallelisms in SOR

Examining equation 12, we note that the values of  $U_{i-1,j}^{(\ell+1)}$  and  $U_{i,j-1}^{(\ell+1)}$  are needed to evaluate  $U_{i,j}^{(\ell+1)}$ . Now consider the following mesh:

0	0	0	0	0	0
0	1	2	3	4	0
0	2	3	4	5	0
0	3	4	5	6	0
0	4	5	6	7	0
0	0	0	0	0	0

---

\* Whenever  $U_{i,j}$  is a boundary point  $U_{i,j}^{(\ell)} = U_{i,j}^{(\ell+1)}$  for all  $\ell$ .

0 indicates a boundary point. The interior mesh points are divided into seven sets (1,2,3..., 7) where set  $i$  contains all points labeled by  $i$ . To improve any point in set  $i$ , we must first improve all the points in the union of all sets  $j$  such that  $j < i$ .

In a computer with 4 PEs, one iteration could be completed in 7 time steps where a time step is the amount of time to improve a mesh point in one PE by improving one set in each time step. But the computer could improve 28 points in 7 time steps. To improve the efficiency of the method, we look at what could be done in each time step assuming we had done everything possible up to that time step.

Time Step	1	2	3	4	5	6	7	8	9	10	11
What	$1^1$	$2^1$	$3^1$	$4^1$	$5^1$	$6^1$	$7^1$				
Can Be			$1^2$	$2^2$	$3^2$	$4^2$	$5^2$	$6^2$	$7^2$		
Done At					$1^3$	$2^3$	$3^3$	$4^3$	$5^3$	$6^3$	$7^3$
That Time							$1^4$	$2^4$	$3^4$	$4^4$	$5^4$
									$1^5$	$2^5$	$3^5$
											$1^6$

Table 1. Parallelism of SOR

$i^l$  denotes the  $l^{\text{th}}$  improvement of the elements in set  $i$ .

In Table 1 we see that steps 1 to 5 improve less than 8 elements but 8 elements are improved at each time step following step 5. Thus if we

have a computer with 8 PEs we can compute  $U^{(\ell)}$  in  $5 + 2\ell$  time steps. If we assume that our initial guess is  $1^3, 2^2, 3^2, 4^1, 5^1, 6^0, 7^0$ , instead of  $1^0, 2^0, 3^0, 4^0, 5^0, 6^0, 7^0$  then we can start at step 6 and do every iteration in 2 time steps.

This can be generalized for an arbitrary mesh as follows: First we group all the interior mesh points into two sets,

$$\text{ODD} = \{U_{i,j} \mid i + j \text{ is odd}\}, \quad \text{EVEN} = \{U_{i,j} \mid i + j \text{ is even}\}$$

Time step 1 improves all the points in EVEN. Time Step 2 improves all the points in ODD. We will call this method modified successive over-relaxation (MSOR).

Remember all the points in ODD can be calculated using only points from EVEN and all the points in EVEN can be calculated using only points from ODD. Thus we can write a two part algorithm for MSOR [Young, Page 271].

$$\begin{aligned} U_{i,j}^{(\ell+1)} = & \omega_{El} (v (U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell)}) + h (U_{i-1,j}^{(\ell)} + U_{i+1,j}^{(\ell)}) - g \Psi_{i,j}) \\ & + (1-\omega_{El}) U_{i,j}^{(\ell)} \end{aligned} \quad (18.1)$$

for all the points in EVEN

$$\begin{aligned} U_{i,j}^{(\ell+1)} = & \omega_{Dl} (v (U_{i,j+1}^{(\ell+1)} + U_{i,j-1}^{(\ell+1)}) + h (U_{i,j+1}^{(\ell+1)} + U_{i,j-1}^{(\ell+1)}) - g \Psi_{i,j}) \\ & + (1-\omega_{Dl}) U_{i,j}^{(\ell)} \end{aligned} \quad (18.2)$$

for all the points in ODD.

In Young's discussion of MSOR, the optimal  $\omega_{E\ell}$  and  $\omega_{D\ell}$  for the  $\ell^{\text{th}}$  iteration are calculated. We will limit our study to MSOR where  $\omega_{E\ell} = \omega_{D\ell} = \omega_b$ .

Intuitively one would think MSOR and SOR have very close rates of convergence. In fact, their "asymptotic rates of convergence" are equivalent.\* If  $A$  is a convergent  $n \times n$  complex matrix, for all  $\ell$  sufficiently large, the average rate of convergence for  $\ell$  iterations,  $R(A^\ell)$ , is  $\lim_{\ell \rightarrow \infty} R(A^\ell) = -\ln \rho(A) \equiv R_\infty(A)$  [Varga, Page 67]. Thus for a sufficiently large  $\ell$ , the asymptotic rate of convergence equals the average rate of convergence. We know if  $P$  is a permutation matrix that  $\rho(PAP^{-1}) = \rho(A)$  [Birkhoff and MacLane, Page 249]. By multiplying (9) by  $P$  we get

$$PU^{(\ell+1)} = P(D - \omega E)^{-1} \{\omega F + (1 - \omega)D\} P^{-1} U^{(\ell)} \quad (19)$$

The asymptotic rate of convergence for (19) equals

$$-\ln \rho(P(D - \omega E)^{-1} \{\omega F + (1 - \omega)D\} P^{-1}) = -\ln \rho((D - \omega E)^{-1} \{\omega F + (1 - \omega)D\}) \quad (20)$$

which equals the asymptotic rate of convergence for SOR.

This not only shows that SOR and MSOR have the same asymptotic convergence rate, but that SOR applied to any permutation of  $U_I$  gives the same asymptotic rate

---

\* The negative of the logarithm of the spectral radius of a convergent matrix  $A$  is the asymptotic rate of convergence,  $R_\infty(A)$ , for the matrix  $A$  [Varga, Page 67].

of convergence. Thus we can choose the ordering of  $U_I$  which best fits the machine being used.

### 2.3.3 SOR using Straight Storage

Now we shall examine the implementation of SOR on ILLIAC IV [Rudsinski]. We will compare two storage schemes and examine their effect on machine efficiency and their time per execution.\*

The first storage scheme, Straight Storage, stores all the elements of row  $i$  before any of the elements of row  $j$  if  $i < j$ . The elements of each row appear in the order they are found in the mesh. The first element of each row is stored in PE 0 and the last element of the row is stored in PE  $I$  where  $I$  equals  $(n-1) \bmod 64$ ,  $n$  is the number of columns. Figure 2 should help clarify this storage scheme.

How can we implement SOR using Straight Storage? We observe that the PEs which contain the odd elements of mesh row  $I$  contain the even elements of mesh row  $I + 1$ . This enables us to improve the odd elements of Rows  $I$  and  $I + 1$  simultaneously by using a PE integer index which accesses an element

---

\* The timing method is explained in Appendix A. The terms used in timing are defined there also.



	PEM 0	PEM 1	...	PEM I	...	PEM 63
	$\vdots$	$\vdots$		$\vdots$		$\vdots$
Row $U_j$	$U_{j,1}$	$U_{j,2}$		$U_{j,I+1}$		$U_{j,64}$
Row $U_{j+1}$	$U_{j,65}$	$U_{j,66}$		$U_{j,65+I}$		$U_{j,128}$
$\vdots$	$\vdots$	$\vdots$		$\vdots$		$\vdots$
Row $U_{j+K-1}$	$U_{j, n-I}$	$U_{j,n+1-I}$		$U_{j,n}$		--
Row $U_{j+K}$	$U_{j+1,1}$	$U_{j+1,2}$		$U_{j+1,I+1}$		$U_{j+1,64}$
	$\vdots$	$\vdots$		$\vdots$		$\vdots$

Figure 2. Straight Storage of Row  $j$  of  $U$  \*

in row  $I$  ( $I+1$ ) in the even (odd) PEs when  $I$  is even. In a like manner we can access all the even elements of two consecutive rows. If we have an odd number of rows we can improve the odd elements of the last row and the even elements of the first row simultaneously using a similar technique.

Given a mesh with  $m$  rows and  $n$  columns, we use the index described above so that we can improve two rows simultaneously. Each pair of rows can

---

\*  $K$  stands for the number of rows of PEM required to store  $n$  words and  $I = (n-1) \bmod 64$ .

be divided into  $K$  groups where  $K = \lfloor n/64 \rfloor^*$  if  $n \bmod 64 = 0$  otherwise  $K = \lfloor n/64 \rfloor + 1$ . Using Phase 1 we can improve 64 elements at a time  $\lfloor (n-2)/64 \rfloor$  times and then use Phase 2 to improve the remaining interior points.\*\* If we use Phase 1 to improve the first 64 odd (even) interior points of a pair of lines then we must take into account that only 63 of them are found in the same pair of rows of PEM because of the boundary points. This means we must use special indices to reach the 64th odd (even) interior point. This requires no extra computer time because the number of times register  $X$  is changed is not affected. SOR requires not only the point to be improved but its four neighbors and the appropriate value of  $\Psi$ .

For example: To improve  $U_{i,j}$  we need  $U_{i-1,j}$ ,  $U_{i+1,j}$ ,  $U_{i,j+1}$ ,  $U_{i,j-1}$ ,  $U_{i,j}$  and  $\Psi_{i,j}$ .

These values will be accessed by a combination of ACAR and Register  $X$  indexing. To do the indexing,  $K$  will be combined with the CU integer  $CI$ -equal to the first row of PEM containing mesh points to be improved during this execution - and three PE integers.

$PI=(1,K,0,K,0,\dots,0,K)$ ,  $PR=(K,0,K,0,\dots,K,0)$  and  $PL=(K+1,1,K,0,K,0,\dots,K,0)$  if we are improving the odd points;

$PI=(K,1,0,K,0,K,\dots,K,0)$ ,  $PR=(0,K,0,K,\dots,0,K)$  and  $PL=(1,K+1,0,K,0,K,\dots,0,K)$  if we are improving the even points;

---

\*  $\lfloor p \rfloor$  equals  $k$  where  $p$  is a real number,  $k$  is an integer, and  $k \leq p < k+1$ .

\*\* Phase 1 and Phase 2 are defined in Appendix D.

INITIAL CONDITIONS			
Register		Contents	
\$ X		PI	
\$ D0		H	
\$ D1		V	
\$ D2		W	
Register		Contents	
\$ D3		WI	
\$ C0		CI	
\$ C1		CM	
\$ C2		CP	
STEP NO.	ASSIGNMENT STATEMENTS	OPERATIONS AND OVERLAP	TIME IN CLOCKS
1	\$A: = U[\$X] (0)	PE fetch	7
2	\$C3: = \$D3; \$A: = \$A*\$C3	CU fetch Real multiplication	10
3	\$S: = \$A - $\psi$ [\$X] (0)	Real subtraction (PE fetch overlapped by multiplication)	7
4	\$A: = U[\$X] (2)	(PE fetch overlapped by addition)	0
5	\$A: = \$A + U[\$X] (1)	Real addition and PE fetch	14
6	\$X: = PL	(PE fetch overlapped by addition)	0
7	\$C3: = \$D0; \$A: = \$A*\$C3	CU fetch Real multiplication	10
8	\$R: = U[\$X] (0)	(PE fetch overlapped by multiplication)	0
9	\$A: = \$A + \$S	Real addition	7
10	\$X: = PR	(PE fetch overlapped by addition)	0
11	\$B: = RTL (1,, \$R)	Route	3
12	\$R: = U[\$X] (0)	(PE fetch partially overlapped by Route)	4
13	\$S: = \$A	Load from PE Register	1
14	\$A: = RTR (1,, \$R)	Route	3
15	\$A: = \$A + \$B	Real addition	7
16	\$C3: = \$D1; \$A: = \$A*\$C3	(CU fetch overlapped by addition) Real multiplication	9
17	\$X: = PI	(PE fetch overlapped by multiplication)	0
18	\$A: = \$A + \$S	Real addition	7
19	\$C3: = \$D2; \$A: = \$A*\$C3	(CU fetch overlapped by addition) Real multiplication	9
20	U[\$X] (0): = \$A	PE store	7
TOTAL TIME			105

Table 2. Timing of an Execution of Phase 1 SOR using Straight Storage

Let  $W = \omega g$ ,  $V = v/g$ ,  $H = h/g$ , and  $Wl = \frac{1-\omega}{W}$  where  $g$ ,  $h$ , and  $v$  are defined in 1.2.

Now (17) can be written in GLYPNIR as follows:

CP: = C + K;

LOOP CI: = C, K + K, CK DO BEGIN

CM: = CI - K;

CP: = CI + K;

U [PI + CI]: = W\*(V\*(RTR(1,,U[PR+CI])+ RTL(1,,U[PL+CI])) + H\*(U[PI+CM]  
+ U[PI+CP]) -  $\psi$ [PI+CI] + Wl\*U[PI+CI] );

(21)

END;

By placing CI in ACAR0, CM in ACAR1, CP in ACAR2, H in \$D0, V in \$D1, W in \$D2, Wl in \$D3 and PI in \$X (21) could be translated into ASK as outlined in Table 2.\*

#### 2.3.4 SOR using Odd-Even Storage

Now we will look at another storage scheme to see if we can improve the data access time. Given an  $m$  by  $n$  mesh divide each row into  $k + 1$  segments such that  $k$  segments are 128 mesh points long and the last segment is  $\ell$  points long, where  $0 \leq \ell < 128$ . Use two adjacent lines of PEM to store each segment. Store all the red (black) points\*\* of the segment in the first (second) line of PEM. The segments appear in the same relative order as they appeared in Straight Storage. This storage scheme will be called Odd-Even Storage. For an example see Figure 3.

---

\* Appendix A explains the notation and timing method used in Table 4.

\*\* A point that appears in an odd(even) numbered column is considered a red (black) point.

If  $\ell \leq 64$ , then Straight Storage requires  $m \left\lceil \frac{n}{64} \right\rceil^*$  lines of PEM while Odd-Even Storage required  $m \left( \left\lceil \frac{n}{64} \right\rceil + 1 \right)$  lines of PEM. If  $\ell > 64$  then both Straight Storage and Odd-Even Storage require  $m \left\lceil \frac{n}{64} \right\rceil$  lines of PEM.

	PEM 0	PEM 1	...	PEM I	...	PEM 63
	$\vdots$	$\vdots$		$\vdots$		$\vdots$
Row $U_j$	$U_{j,1}$	$U_{j,3}$		$U_{j,1+2I}$		$U_{j,127}$
Row $U_{j+1}$	$U_{j,2}$	$U_{j,4}$		$U_{j,2I+2}$		$U_{j,128}$
	$\vdots$	$\vdots$		$\vdots$		$\vdots$
Row $U_{j+K-2}$	$U_{j,n-\ell+1}$	$U_{j,n-\ell+3}$		$U_{j,n-1}$		---
Row $U_{j+K-1}$	$U_{j,n-\ell+2}$	$U_{j,n-\ell+4}$		$U_{j,n}$		---
	$\vdots$	$\vdots$		$\vdots$		$\vdots$

Figure 3. Odd-Even Storage Data Allocation of Row  $j$  of  $U$  Where  $n$  is Even.

---

\*  $\lceil p \rceil = k$  where  $k$  is an integer such that  $p \leq k < p + 1$ .

INITIAL CONDITIONS					
Register		Contents		Register	Contents
\$ X		PI		\$ D3	W1
\$ DO		H		\$ CO	CI
\$ D1		V		\$ C1	CM
\$ D2		W		\$ C2	CP

STEP NO.	ASSIGNMENT STATEMENTS	OPERATIONS AND OVERLAP	TIME IN CLOCKS
1	\$A: = U[\$X] (0)	PE fetch	7
2	\$C3: = \$D3; \$A: = \$A*\$C3	CU fetch Real multiplication	10
3	\$S: = \$A - ψ[\$X] (0)	(PE fetch overlapped with mult.) Real subtraction	7
4	\$C3: = \$CO + 1	(CU operations overlapped)	0
5	\$R: = RTL (1,,U(3))	(PE fetch overlapped by addition) Route	3
6	\$A: = \$R + U[\$X](3)	(PE fetch partially overlapped by Route) Real addition	11
7	\$C3: = \$D1; \$A: = \$A*\$C3	(CU fetch overlapped by addition) Real multiplication	9
8	\$R: = U[\$X](1)	(PE fetch overlapped by multiplication)	0
9	\$S: = \$A + \$S	Real addition	7
10	\$A: = \$U[\$X] (2) + \$R	(PE fetch overlapped by addition) Real addition	7
11	\$C3: = \$DO; \$A: = \$A*\$C3	(CU fetch overlapped by addition) Real multiplication	9
12	\$A: = \$A + \$S	Real addition	7
13	\$C3: = \$D2; \$A: = \$A*\$C3	(CU fetch overlapped by addition) Real multiplication	9
14	U[\$X] (0): = \$A	PE store	7
TOTAL TIME			93

Table 3. Timing of an Execution of Phase 1 SOR using Odd-Even Storage

Phase 1 of the implementation of SOR using Odd-Even Storage requires only one PE index and one route. We can access the data using  $PI = (2,0,0,0..,0)$  and  $CI$  equal the first row of PEM containing mesh points to be improved during this execution. We need two assignment statements: (22.1) for the red points, and (22.2) for the black points.

Using  $K$ ,  $W$ ,  $V$ ,  $H$  and  $Wl$  are defined in 2.2.3, SOR using Odd-Even Storage can be written in GLYPNIR as follows:

```
LOOP CI: = C, K, CK DO BEGIN
```

```
CM: = CI - K;
```

```
CP: = CI + K;
```

```
U[PI + CI]: = W*(V*(RTR(1,, U[CI+1]) + U[PI + CI + 1])
+ H*(U[PI + CM] + U[PI + CP]) -  $\Psi$ [PI + CI] +  $Wl$ *U[PI + CI]); (22.1)
```

```
U[CI]: = W*(V*(RTL(1,,U[PI + CI - 1]) + U[CI - 1])
+ H*(U[CM] + U[CP]) -  $\Psi$ [CI] +  $Wl$ *V[CI]); (22.2)
```

```
END;
```

Table 3 suggests how (22.1) could be coded in ASK. (22.2) requires different code but the logic is similar. The code for SOR using Odd-Even Storage is 10% faster than the code for SOR using Straight Storage. This is due to a savings in memory fetching and the elimination of one route.

### 2.3.5 SOR Machine Efficiency

If a mesh has  $m$  rows and 32 (64) columns and straight (odd-even) storage is used, machine efficiency can be doubled by placing rows 1 to  $\frac{m}{2}$  in PE's 0 to 31 and rows  $\frac{m}{2} + 1$  to  $m$  in PE's 31 to 63. This improved storage scheme requires that rows  $\frac{m}{2}$  and  $\frac{m}{2} + 1$  be handled as special cases. The idea behind this technique can be used on any mesh which doesn't have a multiple



of 64 (128) columns when straight (odd-even) storage is used. Implementation of such techniques increases machine efficiency and decreases storage requirements per mesh, but also increases the complexity of the program.

## 2.4 Implementation of SLOR on ILLIAC IV

### 2.4.1 Parallelism in SLOR

In 2.3.2 the scheme for SOR was altered to maximize parallelism on ILLIAC IV. In a similar manner, we will modify the scheme for SLOR to maximize the number of columns which can be improved simultaneously. In 2.4.2 the need for blocks of 128 columns to perform column relaxation with optimal efficiency on ILLIAC IV will be explained. This leaves no restrictions on the number of rows other than the limits of available core. Analogous to the SOR scheme, note that column  $i$  can be improved for the second time after column  $i + 1$  is improved for the first time. By taking advantage of this fact we can improve half of the columns simultaneously after a finite number of iterations. Thus once we reach that state we can improve the entire mesh in two steps if we have enough PEs. In SOR we improve the odd (even) points and then the even (odd) points. In SLOR we must improve the odd (even) columns and then the even (odd) columns. The proof we used in 2.1.2 to show that we could improve the points in SOR in any order can be extended to show that the order in which the columns are improved SLOR does not change the asymptotic rate of convergence.

$$U_{i,j}^{(\ell+1)} = h(U_{i+1,j}^{(\ell+1)} + U_{i-1,j}^{(\ell+1)}) + \omega v(U_{i,j+1}^{(\ell)} + U_{i,j-1}^{(\ell)}) - \omega g \psi_{i,j}^{(\ell)} + (1 - \omega) (U_{i,j}^{(\ell)} - h(U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)})) \quad (23.1)$$

for the odd columns and

$$U_{i,j}^{(\ell+1)} = h(U_{i+1,j}^{(\ell+1)} + U_{i-1,j}^{(\ell+1)}) + \omega v(U_{i,j+1}^{(\ell+1)} + U_{i,j-1}^{(\ell+1)}) - \omega g \psi_{i,j}^{(\ell+1)} + (1 - \omega) (U_{i,j}^{(\ell)} - h(U_{i+1,j}^{(\ell)} + U_{i-1,j}^{(\ell)})) \quad (23.2)$$

for the even columns.



### 2.4.2 Storage for SLOR

Now we need to find a storage scheme which is efficient on ILLIAC IV. If we use Straight Storage, we run into the same inefficiency we did with that storage in implementing SOR. We need to use three PE integer indices and do two routes per mesh point. Furthermore, the method works on blocks of 256 columns. Column  $J$  and column  $J + 1$  cannot be improved simultaneously. Thus we improve the odd (even) columns in the set  $\{J | I \leq J \leq I + 63\}$  and the even (odd) columns in the set  $\{J | I + 128 \leq J \leq I + 191\}$ . If we use Odd-Even Storage we can eliminate one route and two of the PE indices. We also decrease the blocks to 128 columns each.

### 2.4.3 Improving a Line by the Cuthill-Varga Method

The method of implementing SLOR must solve a tridiagonal matrix equation in each PE. Thomas' method [Todd, Page 395] is commonly used. A method by Cuthill and Varga can eliminate some of the work by doing some preconditioning if the tridiagonal matrix is positive definite and real symmetric. Even if the mesh spacings are not constant, Poisson's equation satisfies those conditions [Cuthill and Varga, Page 241].

We will apply the latter method, the Cuthill-Varga method, as follows. Because we are performing successive column over-relaxation we will use a permutation of  $A$ ,  $PAP^{-1}$ , to discuss the technique. The equation under consideration will be  $-PAP^{-1}PU_I = -PM$ .

$$-PAP^{-1} = \begin{bmatrix} B, C & & & & \\ C, B, C & & & & \\ & C, B, C & & & \\ & & C, B, C & & \\ & & & C, B, C & \\ & & & & C, B, C \\ & & & & & C, B \end{bmatrix} \quad (24)$$

where

$$B = \begin{bmatrix} e, d & & & \\ d, e, d & & & \\ & d, e, d & & \\ & & d, e & \end{bmatrix}, \quad C = \frac{-1}{h_y^2} I, \quad e = +2 \left( \frac{1}{h_y^2} + \frac{1}{h_x^2} \right), \quad d = \frac{-1}{h_x^2}.$$

B and C are  $m-2$  by  $m-2$  matrices. Let  $U_i$  and  $M_i$  denote the  $i^{\text{th}}$  column of PU and -PM respectively. Now using (24) we obtain

$$BU_i = M_i - C(U_{i+1} + U_{i-1}). \quad (25)$$

By assuming the left side of (25) is constant we are left with a tridiagonal system of equations. Because B satisfies the requirements for the Cuthill-Varga method, [Cuthill and Varga, Page 237], we can define a diagonal matrix

$$G = \begin{bmatrix} c_1 & & & \\ & c_2 & & \\ & & \ddots & \\ & & & c_{m-2} \end{bmatrix}$$

such that

$$G^{-1}BG^{-1} = T'T,$$

where<sup>\*</sup>

$$T = \begin{bmatrix} 1 & t_1 & & & \\ & 1 & t_2 & & \\ & & \ddots & \ddots & \\ & & & 1 & t_{m-3} \\ & & & & 1 \end{bmatrix}.$$

The elements of  $G$  and  $T$  are calculated as follows:

$$c_1 = e^{\frac{1}{2}}; \quad c_j = \left\{ e - \left( \frac{d}{c_{j-1}} \right)^2 \right\}^{\frac{1}{2}}, \quad 2 \leq j \leq m-2, \quad (26)$$

and

$$t_j = \frac{d}{c_j c_{j+1}} \quad 1 \leq j \leq m-3. \quad (27)$$

By multiplying

$$BU_i = M_i - C(U_{i+1} + U_{i-1})$$

by  $G^{-1}$  we obtain

$$G^{-1}BG^{-1}GU_i = G^{-1}M_i - G^{-1}CG^{-1}G(U_{i+1} + U_{i-1}).$$

Substituting  $Y_i$  for  $GU_i$  gives

$$G^{-1}BG^{-1}Y_i = G^{-1}M_i - G^{-1}CG^{-1}(Y_{i+1} + Y_{i-1})$$

or

$$T'TY_i = G^{-1}M_i - G^{-1}CG^{-1}(Y_{i+1} + Y_{i-1}) \quad (28)$$

Equation (28) is first solved for  $TY_i$  and then for  $Y_i$ .

The Cuthill-Varga method employs a two part algorithm to perform SLOR.

First  $Y_i^*$  is calculated using

$$T'TY_i^{*(\ell+1)} = G^{-1}M_i - G^{-1}CG^{-1}(Y_{i+1}^{(\ell)} + Y_{i-1}^{(\ell)}) \quad (29.1)$$

if  $i$  is odd and

---

\*  $T'$  equals the transpose of  $T$ .

$$T'_{TY_i}^{*(\ell+1)} = G^{-1}M_i - G^{-1}CG^{-1} (Y_{i+1}^{(\ell+1)} + Y_{i-1}^{(\ell+1)}) \quad (29.2)$$

if  $i$  is even.

Then  $Y_i^{*(\ell+1)}$  is combined with  $Y_i^{(\ell)}$  to perform the over-relaxation,

$$Y_i^{(\ell+1)} = \omega [Y_i^{*(\ell+1)} - Y_i^{(\ell)}] + Y_i^{(\ell)} \quad (30)$$

After all the iterations are completed,  $Y_i$  is converted to  $U_i$  using

$$G^{-1}Y_i = U_i.$$

#### 2.4.4 SLOR Implementation

Now let us discuss more specifically how we implement SLOR\* to solve  $\frac{\delta U}{\delta X^2} + \frac{\delta U}{\delta Y^2} = \Psi$ . Odd-Even Storage issued for the reasons stated in Section 2.4.2. We use the Cuthill-Varga method and thus do the following preconditioning: calculate  $c_i$ ,  $1 \leq i \leq m-2$ ;  $t_i$ ,  $1 \leq i \leq m-3$ ;  $\frac{1}{c_i}$ ,  $1 \leq i \leq m-2$ ; and  $\frac{1}{(h_y c_i)^2}$ ,  $1 \leq i \leq m-2$ . We use PEM storage for these values so that when we calculate  $c_i$  we can do 64 of the square roots simultaneously because the square root takes a great deal of time relative to other operations on ILLIAC IV. We use Grabone\*\* to access the  $c_i$ ,  $t_i$ ,  $\frac{1}{c_i}$  and  $\frac{-1}{(h_y c_i)^2}$ . We subtract the row boundary conditions times  $\frac{1}{h_x^2}$  to the appropriate row of  $\Psi$  and then we multiply by  $G^{-1}$ . We multiply column boundary conditions by  $G$  to obtain  $Y_1$  and  $Y_n$  so that they may be used in equations (29) and (30).

Now we are ready to do the iterations.\*\*\* Two phases are used. Phase 1 improves 64 columns simultaneously and Phase 2 improves the remainder. Phase 1 is used when the number of interior columns is greater than or equal to 128. Phase 1 starts at the left and works on groups of 128 interior columns

---

\* We limit our program to handle an even number of columns.

\*\* The GLYPNIR function is used to access a single PE value and send it to all PEs. Refer to Table 9.

\*\*\* The reader should be familiar with the material in Appendix A.

moving to the right until the number of remaining interior columns is less than 128. Since there are elements of 63 interior odd columns on the first line of PEM containing interior points we use PI and CI as defined in 2.3.4. m-2 rows of temporary storage, STORE, are used. We use two similar codes for Phase 1: one for the odd columns and one for the even columns. Phase 2 can use the same code as Phase 1 if the mode is changed at the appropriate time.

We will limit our discussion to the code for the odd columns, Phase 1. The rest of the code can be found in Appendix D.\*

```
STORE[0] :=  $\psi$ [PI+CI]-GRABONE(DISB[0],0)*(RTR(1,,U[CI+1])+U[PI+CI+1]);
LOOP CJ:= 1,1,CM3 DO BEGIN C:= CJ.[16:42];
CI:= CI+K;
STORE [CJ] :=  $\psi$ [PI+CI]-GRABONE(DISB[C],CJ)*(RTR(1,,U[CI+1])+
U[PI+CI+1])-GRABONE(E[C],CJ)*STORE[CJ-1];
END;
```

(31)

```
LOOP MC:=1,1,CM3 DO BEGIN
CJ:=CM3-MC;
C:=CJ.[16:42];
STORE[CJ] := STORE[CJ]-GRABONE(E[C],CJ)*STORE[CJ+1];
END;
```

(32)

---

\* DISB contains  $\frac{-1}{(h_y c_i)^2}$ ,  $\psi$  contains  $G^{-1}M$  and E contains  $t_i$ .

INITIAL CONDITIONS FOR FORWARD ELIMINATION				
Register		Contents	Register	Contents
\$A		STORE[CJ-1]	\$CO	CI
\$X		PI	\$C1	CJ
			\$C2	CI+1
STEP NO.	ASSIGNMENT STATEMENTS	OPERATIONS AND OVERLAP		TIME IN CLOCKS
1	\$B: = GRABONE(E[C],CU)	Load		10
2	\$A: = \$A*\$B	Real multiplication		9
3	\$S: = $\psi$ [\$X](O)-\$A	(PE fetch overlapped by mult.)		7
4	\$R: = RTR(1,U(2))	Real subtraction		3
5	\$A: = U[\$X](2)+\$R	Route (PE fetch overlapped by subtract.)		3
		(PE fetch partially overlapped by route)		
		Real addition		11
6	\$R: = GRABONE(DISB[C],CU)	Load		10
7	\$A: = \$A+\$R	Real multiplication		9
8	\$A: = \$S-\$A	Real subtraction		7
9	STORE(1): = \$A	PE store		7
Subtotal				73
INITIAL CONDITIONS FOR BACKWARD ELIMINATION				
Register		Contents	Register	Contents
\$A		STORE[CJ+1]	\$CO	CJ
STEP NO.	ASSIGNMENT STATEMENTS	OPERATIONS AND OVERLAP		TIME IN CLOCKS
1	\$B: = GRABONE(E[C],CJ)	Load		10
2	\$A: = \$A*\$B	Real multiplication		9
3	\$A: = STORE(O)-\$A	(PE fetch overlapped by mult.)		7
4	STORE(O): = \$A	Real subtraction PE store		7
Subtotal				33
INITIAL CONDITIONS FOR OVER-RELAXATION				
Register		Contents	Register	Contents
\$X		PI	\$C1	CU
\$CO		CI	\$C2	W
STEP NO.	ASSIGNMENT STATEMENTS	OPERATIONS AND OVERLAP		TIME IN CLOCKS
1	\$R: = U[\$X](O)	PE fetch		7
2	\$A: = STORE(1)-\$R	PE fetch		
		Real subtraction		14
3	\$A: = \$C2*\$A	(CU fetch overlapped by sub.)		
		Real multiplication		9
4	\$A: = \$A+\$R	Real addition		7
5	U[\$X](O): = \$A	PE store		7
Subtotal				44
TOTAL TIME				150

Table 4. Timing of an Execution of Phase 1 SLOR

```

LOOP CJ:=0,1,CM3 DO BEGIN
ST:=U[PI+CI];
U[PI+CI]:=W*(STORE[CJ]-U[PI+CI])+U[PI+CI];
IF ABS(ST-U[PI+CI]) GRT BOUND THEN B:=1;% CHECKING THE ERROR BOUND (33)
CI:=CI+K;
END;

```

SLOR has three primary sections: (31) does the forward elimination, (32) does the backward elimination, and (33) performs the over-relaxation. In Table 4 we suggest how to code these sections efficiently in ASK and we give a time estimate for that code.

## 2.5 Implementation of ADI on ILLIAC IV

ADI improves all columns simultaneously and then improves all rows simultaneously. Thus, when we do the column (row) improvement we want the values of the boundary columns (rows), but the boundary rows (columns) could have been added to vector  $\psi$  in preconditioning. One solution is to include both the row and the column boundary values in each iteration. Another solution is to add them both in when preconditioning and then set the boundary values to zero so they will not affect the interior points when used in the iteration. If we need the values of the boundary after the problem has converged then we must store them before performing any of the iterations and replace them after all the iterations are completed.



### 2.5.1 Thomas' Method on ILLIAC IV

As in SLOR we have sets of tridiagonal matrices to solve. We used the Cuthill-Varga method in SLOR. This method does preconditioning to cut down on the computational time (see Section 2.4.3). ADI changes the diagonal at every iteration. Thus the preconditioning would have to be done at each iteration. Thomas' Method [Todd, page 395] has the same problem but the preconditioning requires less computer time and the total computer time per iteration is smaller. Thus we will use Thomas' Method in solving the tridiagonal matrix problems. Instead of implementing the method in a straight forward manner, we use the fact that the horizontal coefficients and the vertical coefficients are constant.

The matrix problem\* we want to solve is of the form

$$aT_1 + bT_2 = D_1$$

$$bT_{i-1} + aT_i + bT_{i+1} = D_i \quad (i = 2, 3, \dots, n-3),$$

$$bT_{n-3} + aT_{n-2} = D_{n-2}$$

where

$$a = -\left(\frac{2}{h_y^2} + \omega_{\ell V}\right), \quad b = \frac{1}{h_y^2} \quad \text{for row relaxation}$$

and  $a = -\left(\frac{2}{h_x^2} + \omega_{\ell H}\right), \quad b = \frac{1}{h_x^2} \quad \text{for column relaxation}$

Thomas' method is as follows

$$e_1 = \frac{b}{a}, \quad e_i = \frac{b}{a-b e_{i-1}} \quad (i = 2, 3, \dots, n-3), \quad (34)$$

$$q_1 = \frac{D_1}{a}, \quad q_i = \frac{D_i - b q_{i-1}}{a-b e_{i-1}} \quad (i = 2, 3, \dots, n-2), \quad (35)$$

$$T_{n-2} = q_{n-2}, \quad T_i = q_i - e_i T_{i+1} \quad (i = n-3; n-4, \dots, 1). \quad (36)$$

---

\* We are considering an  $m$  by  $n$  mesh,  $U$ .



If we let  $c = \frac{1}{b}$  we can rewrite the equation (35) as follows:

$$q_1 = e_1 c D_1, \quad q_i = e_i (c D_i - q_{i-1}) \quad (i = 2, 3, \dots, n-2),$$

and  $e_{n-2} = \frac{b}{a-b e_{n-3}}$ . This reduces the number of divisions which are relatively time consuming on ILLIAC IV.

We have to calculate the  $e_i$ 's and store them. For each row (column), they are the same. Calculating the  $e_i$ 's is a serial process. The only way we could overlap is to calculate the set of  $e_i$ 's for different iterations simultaneously. The problem with this solution is that it requires extra storage. If we calculate all the  $e_i$ 's needed for 32 iterations we use  $\max(m-2, n-2)$  rows of storage. If storage is available, then this is a feasible method; but if we must use disk I/O, the computational time saved will be overcome by increased I/O time.

### 2.5.2 Skewed Storage

The next problem we must overcome is how to access the rows on one sweep and then the columns on the next. Skewed storage was designed just for that purpose. We can place a  $m$  by  $n$  mesh in skewed Storage by performing the following algorithm:

- 1) Calculate the number of rows of PEM required to store  $n$  points. Set  $K$  equal to that number.
- 2) Place the  $m$  by  $n$  mesh in straight storage.
- 3) Route each of the  $K$  rows of PEM containing elements of mesh row  $i$ . Let the distance of the route be  $i-1$ .

	PEM 0	PEM 1	...	PEM 31	PEM 32	PEM 33	...	PEM 62	PEM 63
	:	:		:	:	:		:	:
Row $U_1$	$U_{1,1}$	$U_{1,2}$		$U_{1,32}$	$U_{1,33}$	$U_{1,34}$		$U_{1,63}$	$U_{1,64}$
Row $U_{1+1}$	$U_{1,65}$	$U_{1,66}$		$U_{1,96}$	--	--		--	--
Row $U_2$	$U_{2,64}$	$U_{2,1}$		$U_{2,31}$	$U_{2,32}$	$U_{2,33}$		$U_{2,62}$	$U_{2,63}$
Row $U_{2+1}$	--	$U_{2,65}$		$U_{2,95}$	$U_{2,96}$	--		--	--
:	:	:		:	:	:		:	:
Row $U_{33}$	$U_{33,33}$	$U_{33,34}$		$U_{33,64}$	$U_{33,1}$	$U_{33,2}$		$U_{33,31}$	$U_{33,32}$
Row $U_{33+1}$	--	--		--	$U_{33,65}$	$U_{33,66}$		$U_{33,95}$	$U_{33,96}$
Row $U_{34}$	$U_{34,32}$	$U_{34,33}$		$U_{34,63}$	$U_{34,64}$	$U_{34,1}$		$U_{34,32}$	$U_{34,31}$
Row $U_{34+1}$	$U_{34,96}$	--		--	--	$U_{34,65}$		$U_{34,99}$	$U_{34,95}$
:	:	:		:	:	:		:	:
Row $U_{96}$	$U_{96,34}$	$U_{96,35}$		$U_{96,1}$	$U_{96,2}$	$U_{96,3}$		$U_{96,32}$	$U_{96,33}$
Row $U_{96+1}$	--	--		$U_{96,65}$	$U_{96,66}$	$U_{96,67}$		$U_{96,96}$	--
		:		:	:	:		:	:

Figure 4. A 96 by 96 Mesh in Skewed Storage

ADI is a simultaneous method, (i.e. only values from the previous iteration are used, see (15.1) and (15.2)). On ILLIAC IV we can simultaneously improve up to 64 rows (columns). Thus if the dimensions of the mesh are less than 67 we can do all the rows (columns) in one time step. If one of the dimensions is greater than 66 then we need to use more than one time step for row or column relaxation. In this case we will use two phases to complete an iteration. An  $m$  by  $n$  mesh will be divided using  $R = \left\lfloor \frac{m-2}{64} \right\rfloor$  and  $J = \left\lfloor \frac{n-2}{64} \right\rfloor$ . Phase 1 row (column) relaxation will improve  $R$  ( $J$ ) groups of 64 consecutive rows (columns). Then Phase 2 will improve the remaining interior rows (columns). Phase 1 starts with row (column) 2 and works toward row (column)  $m-1$  ( $n-1$ ). When Phase 1 improves the  $I^{\text{th}}$  group where  $1 \leq I \leq R$  ( $J$ ) and  $k$  is the first row (column) of group  $I$ , the old values of row (column)  $k-1$  are needed. Thus to improve group  $I$  Phase 1 goes through the following steps:

- 1) The values of row (column)  $k-1$  are placed in the temporary storage, NEW.
- 2) The values in the temporary storage, OLD are placed in the PEM storage for row (column)  $k-1$ .
- 3) The values of row (column)  $k+63$  are placed in the temporary storage, OLD.
- 4) Rows (columns)  $k$  through  $k+63$  are improved.
- 5) The values in NEW are placed in row (column)  $k-1$ .

When Phase 1 is improving the first group steps 1, 2 and 5 can be omitted. Phase 2 needs to do steps 1, 2, 4 and 5.

Now we will discuss the type of indexing used in accessing the mesh points required to perform Phase I column relaxation using skewed storage. We access\* the set  $\{(\psi_{i,j}, U_{i,j}) | k \leq j \leq k + 63\}$  by using a permutation of the PE integer,  $(1, 0, 0, \dots, 0)$ . PI will stand for that permutation.

PI will be used as a PE index and CI will be the CU index. PI must be routed one PE to the right to access the set  $\{U_{i+1,j} | k \leq j \leq k+63\}$ . PI must be routed one PE to the left to access  $\{U_{i-1,j} | k \leq j \leq k+63\}$ . To access the set  $\{U_{i,j+1} | k \leq j \leq k+63\}$  we must add  $RTR(1, PI)$  to PI. The set  $\{U_{i,j-1} | k \leq j \leq k+63\}$  can be accessed using only CI. After being accessed the four neighbors of  $U_{i,j}$  must be routed to the PE containing  $U_{i,j}$  before  $U_{i,j}$  can be improved. The following GLYPNIR code uses this method of data access to perform Phase 1, Step 4 of column relaxation:

```

LOOP C:=1,1,N-2 DO BEGIN CC:= C. [16:42]; CI:=CI+K;
U[PI+CI]:= (AI*( $\psi$  [PI+CI]-BETA*(RTL(1,,U[PI+RPI+CI])+RTR(1,,U[CI]))
+CA*U[PI+CI])-RTR(1,,U[LPI+CI-K]))*GRABONE(B[C-1],CC);
LPI:=PI; PI:=RPI; RPI:=RTR(1,,RPI); END;

```

```

LOOP IC:=0,1,N-3 DO BEGIN C:=N-2-IC;
CC:=C.[16:42]; CI:=CI-K;
U[PI+CI]:=U[PI+CI]-GRABONE(B[C-1],CC)*RTL(1,,U[RPI+CI+K]);
RPI:=PI; PI:=RTL(1,,PI); END;

```

---

\* The variable, k equals the first column of the group being improved.

\*\*  $AI = -h_y^2$  ;  $BETA = 1/h_x^2$  ; W equals the relaxation parameter for this iteration;  $CA = 2*BETA + W$ .

The code is divided into two sections: Section (37) does Thomas' forward elimination. Section (38) does Thomas' backward elimination; Table 7 suggests how this code could be translated into ASK and how much time the ASK code would take.

By replacing PI with  $KPEN=RTR(I,,PENK)$  where  $PENK=(0,K,2k,\dots,63K)$  and K is the number of lines of PEM required to store a row of the mesh, we would have Phase 1 row relaxation.\*

A complete program for a m by n mesh where  $n, m \leq 64$  is found in Appendix E.

---

\* In row relaxation a PE index must be used to access  $\{U_{i,j-1} | k \leq j \leq k+63\}$ .

INITIAL CONDITIONS FOR FORWARD ELIMINATION					
Register		Contents	Register		Contents
\$C1		CI	\$D2		CA
\$D0		AI	\$X		PI
\$D1		BETA	\$C3		CI-K
STEP NO.	ASSIGNMENT STATEMENT		OPERATIONS AND OVERLAP		TIME IN CLOCKS
1	\$A: = U[\$X](1)		PE fetch		7
2	\$C2: = \$D2;\$A:=\$A*\$C2		CU fetch and real mult.		10
3	\$S: = \$A + $\psi$ [\$X](1)		(PE fetch overlapped by mult.) Real addition		7
4	PI: = \$X		(PE store overlapped by add.)		0
5	\$X: = \$X + RPI		PE fetch and integer addition		10
6	\$B: = RTL(1,,U[\$X](1))		(PE fetch partially overlapped by addition) Route		7
7	\$A: = RTR(1,,U(1))		(PE fetch partially overlapped by route) Route		7
8	\$X: = LPI		(PE fetch partially overlapped		4
9	\$A: = \$A + \$B		Real addition		7
10	\$R: = U[\$X](3)		(PE fetch overlapped by add.)		0
11	\$C2: = \$D1;\$A:=\$A*\$C2		CU fetch, real multiplication		10
12	\$X: = PI		(PE fetch overlapped by mult.)		0
13	\$A: = \$S-\$A		Real subtraction		7
14	LPI: = \$X		(PE store overlapped by subtr.)		0
15	\$C2: = \$D0;\$A:=\$A*\$C2		CU fetch, real multiplication		10
16	\$B: = RTR(1,, \$R)		Route		3
17	\$A: = \$A+\$B		Real addition		7
18	\$S: = RPT		(PE fetch overlapped by add.)		0
19	\$R: = RTR(1,, \$S)		Route		3
20	\$A: = \$A*GRABONE(B[CI-1],CC)		Load real multiplication		19
21	RPI: = \$R		(PE store overlapped by mult.)		0
22	U[\$X](1): = \$A		PE store		7
23	\$X: = \$S		PE register to PE register load		1
Subtotal					126
INITIAL CONDITIONS FOR BACKWARD ELIMINATION					
Register		Contents	Register		Contents
\$X		PI	\$A		U[RPI+CI+K]
			\$C1		CI
STEP NO.	ASSIGNMENT STATEMENT		OPERATIONS AND OVERLAP		TIME IN CLOCKS
1	\$B: = GRABONE(B[C-1],CC)		Load		10
2	\$A: = \$B*\$RTL(1,, \$A)		Route, real multiplication		12
3	\$A: = U[\$X](1)-\$A		(PE fetch overlapped by mult.) Real subtraction		7
4	U[\$X](1): = \$A		PE store		7
5	\$X: = RTL(1,, \$X)		(Route overlapped by store)		0
Subtotal					36
TOTAL TIME					162

Table 5. Timing of an Execution of Phase 1 ADI



## 2.6 Summary of the Results for Iterative Methods

Now that we have examined implementation of the various methods we can make some recommendations concerning the use of the methods. Table 6 contains a numerical study of convergence rates of the method applied to Poisson's equation where  $h_x = h_y = \frac{1}{15}$  on different mesh sizes. The error bound was 0.0001.

	64x64	32x64	32x32	16x64	Mesh size Method
$\omega$	1.86415	1.79300	1.74760	1.64750	SLOR Column Relaxation
$I_t$	63	40	32	21	
$I_a$	66	41	33	23	
$\omega$	1.90645	1.85610	1.82147	1.75050	SOR
$I_t$	92	59	41	32	
$I_a$	90	63	44	28	
$I_a$	10	9	8	7	ADI
$\eta$	9	8	7	6	

- a)  $\omega$  is the relaxation parameter
- b)  $I_t$  is the theoretical number of iterations
- c)  $I_a$  is the actual number of iterations\*
- d)  $2\eta$  is the number of relaxation parameters used in ADI

Table 6. A Numerical Study of the Methods under Discussion

We can see from Table 6 that ADI is consistently the fastest\*\*, then comes SLOR, with SOR being slowest. This conforms with theoretical results. In fact, in a square mesh where  $h_x = h_y$ , SLOR converges  $\sqrt{2}$  times faster than SOR while ADI converges  $\beta$  times faster than SOR.  $\beta$  is a monotonically increasing function of the number of mesh points [Todd, Pages 396-398].

\* Remember ADI improves the points twice in each iteration.

\*\* The speed of the algorithms is compared with respect to the number of iterations.

To compare the time required to perform one of the above methods on ILLIAC IV to that required to perform the method on a specific machine we need to calculate the clocks per point per iteration and compare it with the clocks per point per iteration on ILLIAC IV. We have shown that SOR, SLOR, and ADI can be programmed to enable ILLIAC IV to improve 64 points simultaneously for certain mesh sizes.\* By dividing the clocks required to improve 64 points simultaneously by 64 we obtain the clocks per point per iteration on ILLIAC IV. Table 7 contains these values and other pertinent information.

		ADI	SLOR	SOR
Number of operations per point per iteration	Mult.	10	4	4
	Add.	10	6	5
Time in clocks per 64 points per iteration	Arithmetic	160	78	71
	Data Access	164	72	22
Total clocks per point per iteration		5.06	2.34	1.45
Optimum mesh size**		64I by 64L	128I by L	64I by 2L
Rows of PEM required for temporary storage		Max (n-2, m-2)	$m-2 + \frac{(m-2)}{64}$	0

Table 7. Summary of Time and Storage Requirements for the Iterative Methods on a m by n Mesh.

Most of the data access time on ILLIAC IV could be eliminated if it was a serial machine. Let Machine A be a serial machine with a processor similar in speed to a PE\*\*\* but with the ability to perform ADI, SOR, and SLOR with negligible data access time\*\*\*\*. From Table 7 we can see that Machine A would take  $71/1.45$ , about 48, times longer to perform SOR than ILLIAC IV would take. Machine A would take about 32 times longer than ILLIAC IV to perform ADI or SLOR.

\* For maximum efficiency on ILLIAC IV, one dimension of the mesh must be a multiple of 64 for SOR and SLOR while both dimensions must be multiples of 64 for ADI.

\*\* I and L are positive integers

\*\*\* A single PE has approximately twice the arithmetic speed of a CDC 6600.

\*\*\*\* The clocks per point per iteration on Machine A equal the arithmetic time in clocks per 64 points per iteration on ILLIAC IV.



The times in Table 7 are assuming the inner loops of the programs are written in ASK. If the programs were written completely in GLYPNIR the times would about double.

Tables 6 and 7 indicate that ADI is the fastest iterative algorithm on ILLIAC IV for rectangular meshes. The theory for ADI has only been developed for special cases [Young, Page 555]. If we have a commutative case,  $HV = VH$ , then ADI will be effective. The commutative space requires a rectangular region and coefficients of the differential equation which are sufficiently regular [Young, Page 538]. In some noncommutative cases, numerical experiments have shown ADI to converge rapidly. This is not always true. In some noncommutative cases ADI fails to converge for certain parameters [Young, Page 546]. More work needs to be done on noncommutative cases before ADI can be used freely on them.

ADI performs row relaxation followed by column relaxation. Thus to efficiently perform ADI on a  $m$  by  $n$  mesh on ILLIAC IV, both  $m$  and  $n$  must be divisible by 64. For example if we have a 16 by 64 mesh when we perform row relaxation, ILLIAC IV will be working at 25% machine efficiency while for column relaxation ILLIAC IV would be working at 100% machine efficiency. If one had four 16 by 64 meshes to solve, ILLIAC IV could solve the four meshes simultaneously and thus brings the machine efficiency up to 100%.

SLOR converges in fewer iterations than SOR and can be programmed to take about the same amount of time per iteration as SOR on most serial machines. On ILLIAC IV SLOR requires data from one PE broadcast to all the PEs while SOR does not. This is the major factor in causing SLOR to

take about 1.5 times longer per iteration than SOR, see Table 7. Thus SOR is a faster method on ILLIAC IV unless it takes at least 1.5 times more iterations to converge. This is rarely the case. SLOR requires considerably more temporary storage than SOR. Finally SLOR is a block iterative method and thus cannot be as easily reformulated as SOR for efficient performance on ILLIAC IV.

As shown in Section 2.3 of this study, SOR has a few constraints on how it must be implemented. This enables one to program it efficiently on ILLIAC IV for most mesh geometrics. SOR seems to be the most promising of the iterative and direct methods examined by the author for non-rectangular meshes.

In this study, we indicated that the order in which the elements are improved by SOR does not affect the asymptotic rate of convergence. This can be misunderstood. The number of iterations required to obtain a specified error bound is dependent on the order in which the elements are improved. For specific initial conditions, one ordering might give a faster initial rate of convergence than another. In some applications of Poisson's equation the actual number of iterations required for convergence is substantially smaller than theoretically expected because a good initial guess is supplied. In these cases the ordering has a greater affect on the computer time required to get an acceptable solution. ADI and SLOR require a considerable amount of preconditioning before the iterative process can begin. Thus if the initial guess is good enough, SOR will take the least amount of computer time on any mesh.

This study examined straight and odd-even storage for SOR. The use of odd-even storage improved the speed of SOR by 10%. In most problems, straight storage is used. Thus, for a valid comparison of the two storage schemes, the time to convert straight storage to odd-even storage must be considered. This takes about the same amount of time as one SOR iteration. Thus the program must require at least 20 iterations to justify converting to odd-even storage and then back to straight storage.

### 3. IMPLEMENTATION OF DIRECT METHODS ON ILLIAC IV

#### 3.1 Introduction to Hockney's Direct Method (FACR)

Direct methods for solution of a restricted class of Poisson's equation have been developed which are faster than any iterative method developed to date [Dorr, Pages 258-259]. To the author's knowledge, Hockney's Fourier Analysis/Cyclic Reduction (FACR) is the fastest direct method on serial machines [Hockney, Page 159]. In this paper we will examine FACR and modify it so that it can be programmed efficiently on ILLIAC IV. FACR solves the "five-point" difference formula on a rectangular mesh; namely,

$$\frac{U_{s-1,t} - 2U_{s,t} + U_{s+1,t}}{h_x^2} + \frac{U_{s,t-1} - 2U_{s,t} + U_{s,t+1}}{h_y^2} = \psi_{s,t} \quad (39)$$

where the number of points being computed in one direction is  $2^N - 1$  for Dirichlet's boundary conditions,  $2^N$  for periodic boundary conditions and  $2^N + 1$  for Neumann's boundary conditions. These three boundary conditions are permitted in the x and y direction, giving nine possible combinations of boundary conditions.

A  $2^m$  by  $2^n$  mesh produces a linear system with  $2^{n+m}$  unknowns. Let  $M = 2^m$  and  $N = 2^n$ . FACR solves this system using the following five step algorithm:

1. Given  $\Psi$  compute  $\Psi^*$  for the even columns and overwrite  $\Psi$  with  $\Psi^*$  on the even columns.
2. Using  $\Psi^*$  compute  $\Psi^S$  and overwrite  $\Psi^*$  with  $\Psi^S$ .
3. Using  $\Psi^S$  solve for  $U^S$  and overwrite  $\Psi^S$  with  $U^S$ .
4. Using  $U^S$  compute  $U$  on the even columns and overwrite  $U^S$  with  $U$  on the even columns.

5. Using the values of  $U$  on the even columns solve (39) for the values of  $U$  on the odd columns and overwrite  $\Psi$  with  $U$  on the odd columns.

The superscripts are for Dirichlet's boundary conditions and are defined as follows:

$$\Psi_{i,j}^* = \Psi_{i,j+1} - \frac{h_y^2}{h_x^2} (\Psi_{i-1,j} + \Psi_{i+1,j}) + 2 \left( \frac{h_y^2}{h_x^2} + 1 \right) \Psi_{i,j} + \Psi_{i,j-1}$$

$$\Psi_{i,j}^S = \frac{2}{M} \sum_{k=1}^{M-1} \Psi_{k,j}^* \sin \frac{\pi i k}{M}$$

$$U_{i,j}^S = \frac{2}{M} \sum_{k=1}^{M-1} U_{k,j} \sin \frac{\pi i k}{M}$$

Odd/even reduction divides the problem into two parts; first to solve a linear system concerning even columns and secondly to solve for the values on the odd columns. The even columns form a linear system with  $M \frac{N}{2}$  unknowns which is solved by steps 2, 3, and 4. Step 5 involves the  $\frac{N}{2}$  linear systems with  $M$  unknowns that give the values for the odd columns.

Fourier analysis decouples the even columns into  $M$  tridiagonal linear systems with  $\frac{N}{2}$  unknowns. The unknowns are the Fourier harmonics,  $U^S$ , of  $U$ .

Recursive cyclic reduction solves the linear systems for the Fourier harmonics  $U^S$ .

Fourier synthesis converts the Fourier harmonics  $U^S$  obtained by the previous step to  $U$  giving the values of  $U$  on the even columns.

The values of  $U$  on the even columns are used to calculate the values of  $U$  on the odd columns. This involves solving  $\frac{N}{2}$  tridiagonal linear systems with  $M$  unknowns [Hockney, Pages 148-153].

Refer to Appendix B for a more detailed explanation. Note that all five steps can be performed using one mesh containing the boundary conditions and  $\Psi$  initially. This is overwritten by  $\Psi^*$ ,  $\Psi^S$ ,  $U^S$  and finally the solution  $U$ .

We will discuss the method in three parts, steps 1 and 5, step 3, and steps 2 and 4. Suggestions will be made for the application of each part. Finally we will present a program in GLYPNIR similar to FACR and make suggestions on how it can be programmed for ASK.

### 3.2 Odd/Even Reduction and Odd Column Solution

Odd/even reduction is used to cut down on the number of columns where Fourier analysis and synthesis is applied. Consider the three neighboring equations:

$$\begin{aligned} \frac{1}{h_y^2} U_{t-2} + BU_{t-1} + \frac{1}{h_y^2} U_t &= \Psi_{t-1} \\ \frac{1}{h_y^2} U_{t-1} + BU_t + \frac{1}{h_y^2} U_{t+1} &= \Psi_t \\ \frac{1}{h_y^2} U_t + BU_{t+1} + \frac{1}{h_y^2} U_{t+2} &= \Psi_{t+1} \end{aligned} \quad (40)$$

where  $U_t$  is the  $t^{\text{th}}$  column of the mesh and  $t$  is even. By multiplying the middle even line equation by  $-Bh_y^2$  and adding we obtain

$$\frac{1}{h_y^2} U_{t-2} + \left( \frac{2}{h_y^2} - B^2 h_y^2 \right) U_t + \frac{1}{h_y^2} U_{t+2} = \Psi_{t+1} - Bh_y^2 \Psi_t + \Psi_{t-1} = \Psi_t^* \quad (41)$$

Note equation (41) contains only even columns of  $U$ , that is  $U_{t-2}$ ,  $U_t$  and  $U_{t+2}$ . Thus after (41) is solved the algorithm must solve for the values of  $U$  on the odd columns.



To apply odd/even reduction one needs an odd number of mesh points for Dirichlet or Neumann boundary conditions and an even number of mesh points for periodic boundary conditions.

Odd/even reduction halves the number of columns that require Fourier analysis and synthesis. The odd columns are solved as tridiagonal matrix problems, step 5. Cyclic reduction is used to solve for the odd columns of U. This has the potential of saving computational time. Hockney showed that one level of odd/even reduction reduces the total number of operations per point on a serial machine from  $3^4$  to  $2^4$  on a  $128 \times 128$  mesh [Hockney, Page 161].

Let  $k$  be the number of interior columns in the mesh. On ILLIAC IV Fourier analysis and synthesis can be applied to  $6^4$  of these columns simultaneously. Thus without odd/even reduction a Fourier algorithm must be applied  $\lceil \frac{k}{64} \rceil$  times.\* If odd/even reduction is used a Fourier algorithm must be applied  $\lceil \frac{k}{128} \rceil$  times.

Odd/even reduction is of no use if  $k \leq 64$  because then  $\lceil \frac{k}{64} \rceil = \lceil \frac{k}{128} \rceil$ . If  $k > 64$  then  $\lceil \frac{k}{64} \rceil > \lceil \frac{k}{128} \rceil$  and odd/even reduction would reduce the number of times Fourier analysis and synthesis must be applied. The best case is where  $\lceil \frac{k}{64} \rceil = 2 \lceil \frac{k}{128} \rceil$  in which case the number of Fourier analysis and synthesis is cut in half.

Straight storage is the standard storage scheme used by most programs. If odd/even reduction is used, maximum machine efficiency is obtained when an even column is contained in each PE for the odd/even

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\*  $\lceil P \rceil$  equals the integer  $L$  where  $L - 1 < P \leq L$ .

96 x 96

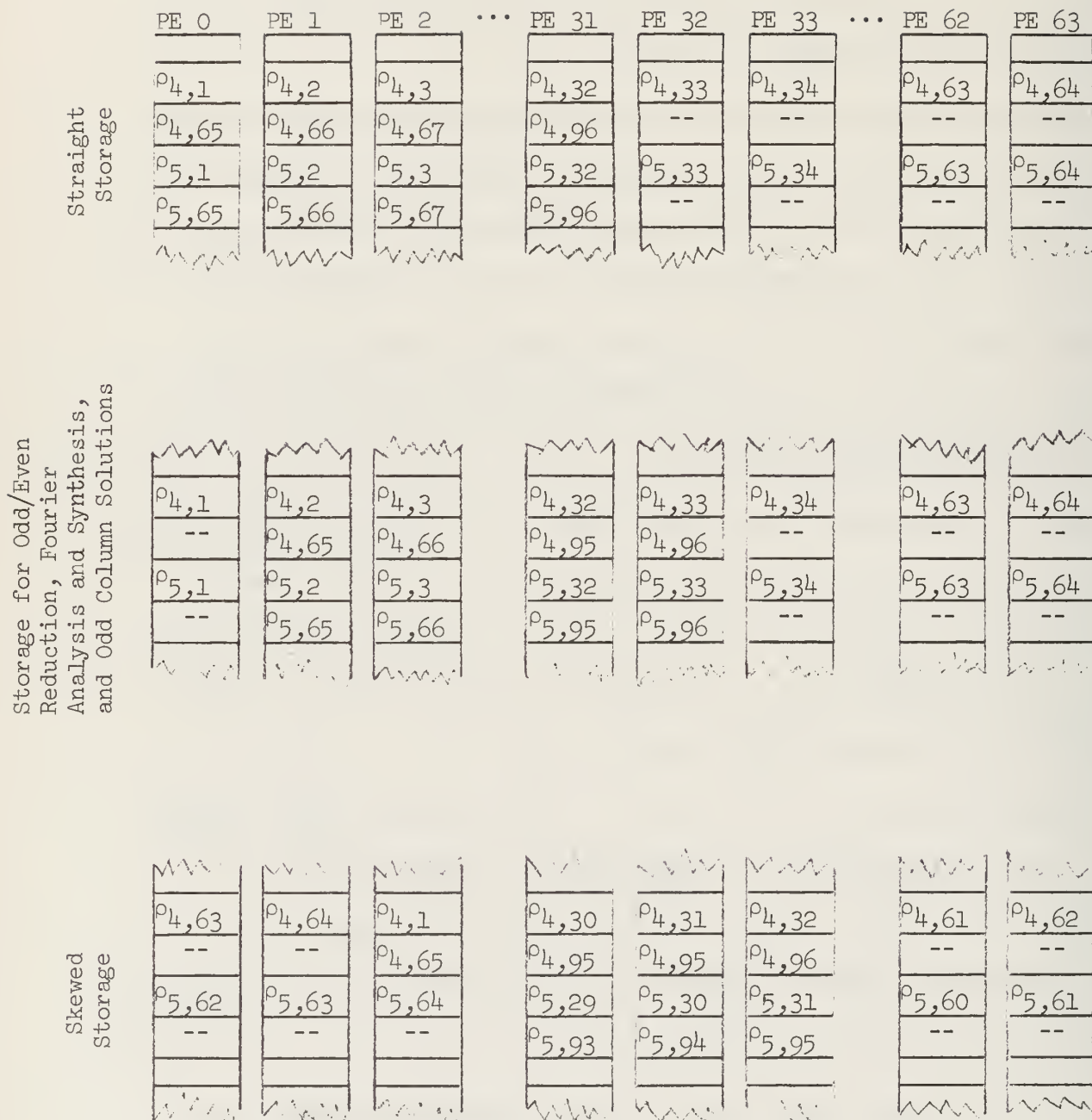


Figure 5. Storage Schemes Used on Rows 4 and 5 of a 96 by 96 Mesh if Odd/Even Reduction is Used



reduction step, the Fourier analysis step and the Fourier synthesis step; CRED can access a different row in each PE; and during the solution of the odd columns an odd column can be accessed by each PE. The following storage changes will fulfill all these conditions assuming each row of the mesh requires an even number of rows of PEM.

1. We assume the main program supplies U in straight storage.
2. Before odd/even reduction, the even rows of PEM containing U are routed one PU to the right of the odd rows.
3. Before CRED, U is placed in skewed storage.
4. Before Fourier synthesis, U is returned to the storage used for odd/even reduction.
5. The mesh U is placed in straight storage at the end of the subroutine.

See Figure 5 for clarification of the above storage schemes. If odd/even reduction is not used steps 2 and 4 of the storage changes need not be executed. The use of odd/even reduction requires the extra storage changes because steps one through four of FACR are applied to even columns while step five solves the odd columns. Without the use of odd/even reduction steps 2, 3, and 4 in Section 3.1 would be applied to all the columns.

### 3.3 CRED

In step 3, section 3.1 we have a tridiagonal system of equations to solve. The derivation of the coefficients for Dirichlet's boundary conditions is contained in Appendix B. By multiplying equation (B-8) in Appendix B by  $h_y^2$  the diagonal coefficient for the  $k^{\text{th}}$  row becomes

$$\lambda_k = - \left( 2 \frac{h_y^4}{h_x^4} \cos \frac{2\pi k}{M} - 8 \left( \frac{h_y^4}{h_x^4} + \frac{h_y^2}{h_x^2} \right) \cos \frac{\pi k}{M} + 6 \frac{h_y^4}{h_x^4} + 8 \frac{h_y^2}{h_x^2} + 2 \right)$$

Hockney solves these tridiagonal systems using cyclic reduction recursively.

Given

$$\begin{aligned} U_{k,t-4}^s + \lambda_k U_{k,t-2}^s + U_{k,t}^s &= h_y^2 \psi_{k,t-1}^s \\ U_{k,t-2}^s + \lambda_k U_{k,t}^s + U_{k,t+2}^s &= h_y^2 \psi_{k,t}^s \quad (42) \\ U_{k,t}^s + \lambda_k U_{k,t+2}^s + U_{k,t+4}^s &= h_y^2 \psi_{k,t+1}^s \end{aligned}$$

where  $t$  is even.

By multiplying the middle equation by  $-\lambda_k$  and adding we obtain

$$U_{k,t-4}^s + (2 - \lambda_k^2) U_{k,t}^s + U_{k,t+4}^s = h_y^2 (\psi_{k,t-2}^s + \psi_{k,t+2}^s - \lambda \psi_{k,t}^s) \quad (43)$$

Now the process is repeated on the equations until we are left with just one equation. Then the process is reversed [Hockney, Page 150]. This method uses  $\log_2 N$  extra words of storage,  $4N - 2\log_2 N$  additions,  $2N$  multiplications and  $\log_2 N$  divisions for an  $N$  variable system.\* It also requires that  $N$  equal  $2^I - 1$  if the system has Dirichlet conditions,  $2^I$  if the system has periodic conditions, and  $2^I + 1$  if the system has Neumann conditions.

For Dirichlet boundary conditions Thomas' method allows one to solve a tridiagonal system of  $N$  variables where there are no restrictions on  $N$ , (See Section 2.5.1). In the case where the diagonal elements are equal and all the other non-zero elements equal 1, we calculate

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\*  $N$  is the number of interior even columns in the mesh.

$$e_{k,1} = \frac{1}{\lambda_k} ; e_{k,i} = \frac{1}{\lambda_k - e_{k,i-1}} \quad \text{for } i = 2, 3, \dots, N. \quad (44)$$

$$q_{k,1} = \psi_{k,1}^s e_{k,1} ; q_{k,i} = (\psi_{k,i}^s - q_{k,i-1}) e_i \quad \text{for } i = 2, 3, \dots, N. \quad (45)$$

$$U_{k,N}^s = q_{k,N} ; U_{k,i}^s = q_{k,i} - e_{k,i} U_{k,i+1}^s \quad \text{for } i = N-1, N-2, \dots, 1. \quad (46)$$

This method uses  $N$  words of extra storage,  $N$  divisions,  $2N$  multiplications, and  $3N$  additions. On ILLIAC IV a division requires as much time as six multiplications. Thus the calculations of the  $e_i$ 's takes most of the time required for Thomas' method. Since the  $e_i$ 's are independent of  $U$  and  $\psi$  they could be calculated in preconditioning and used repeatedly by Thomas' method. Since each row has a different  $\lambda_k$  we need a different set of  $e_{k,i}$ 's for each row of  $U$ . Thus all the  $e_{k,i}$ 's would require half as much storage as  $U$  if odd/even reduction is applied. Thomas' method can be applied to an arbitrary  $M$ . This not only allows the program to be more general but in some cases can save storage in placing  $U$  and  $\psi$  in PEM. If a problem with Dirichlet or Neumann boundary conditions required the accuracy obtained by  $64$  by  $64$  meshes, applying cyclic reduction would require the meshes to be  $65$  by  $65$  while Thomas' method would accept  $65$  by  $64$  meshes. In straight storage a  $65$  by  $65$  mesh requires 130 lines of PEM while a  $65$  by  $64$  mesh requires 65 lines of PEM. If the boundary conditions are periodic, cyclic reduction would require a  $64$  by  $64$  mesh, and would require only 64 lines of PEM. The extra storage required to store the  $e_{k,i}$ 's for Thomas' method is too great for some memory bound problems. By changing (46) to

$$\begin{aligned} U_{k,N}^s &= q_N^s ; U_{k,N-1}^s = \psi_{k,N}^s - \lambda_k U_{k,N}^s ; \\ U_{k,i}^s &= \psi_{k,i+1}^s - \lambda_k U_{k,i+1}^s - U_{k,i+2}^s \quad \text{for } i = N-2, N-3, \dots, 1. \end{aligned} \quad (47)$$

we eliminate the need for extra storage. We will refer to this method as modified Thomas' method. It takes  $N$  divisions,  $2N$  multiplications, and  $4N$  additions.

Table 8 summarizes the storage and computational time of cyclic reduction, Thomas' method and modified Thomas method where the divisions are done in preconditioning for Thomas' method and cyclic reduction and the divisions are not done in preconditioning for the modified Thomas' method. The computational time is calculated by multiplying 7, 9 and 56 clocks by the number of additions, multiplications and divisions respectively and is for  $64$  rows of  $U$ . The storage is in words not lines of PEM.

	Mesh Size Restrictions	Extra Storage Requirements		Time per 64 Rows in Clocks
Thomas' Method with odd/even reduction	An odd number of columns	Thomas' Method	$NM/2$ words	$39N/2$
		Mod. Thomas' Method	No words	$102N/2$
Thomas' Method without odd/even reduction	No Restrictions	Thomas' Method	$NM$ words	$39N$
		Mod. Thomas' Method	No words	$102N$
Cyclic Reduction	$N = 2^I + 1$	$M \log_2 N$ words		$46N - 14 \log_2 N$

Table 8. Comparison of Three Methods to Solve Tridiagonal Matrix Equations for Dirichlet's or Neumann's boundary conditions.

Odd/even reduction is most helpful when  $\left\lceil \frac{k}{64} \right\rceil = 2 \left\lceil \frac{k}{128} \right\rceil$  where  $k$  is the number of interior columns (see Section 2). The time estimates of Table 8 for Thomas' method with odd/even reduction are only half the corresponding values for Thomas' method without odd/even reduction. This is because

odd/even reduction decreases the size of the tridiagonal system to be solved by a factor of 2.

For cyclic reduction the amount of extra storage in Table 8 is misleading. If  $I \geq 6$  then  $64M$  words of extra storage are required per mesh in PEM. If both  $U$  and  $\Psi$  are in PEM, then  $128M$  words are required. This is because cyclic reduction requires  $2^I + 1$  columns for Dirichlet's or Neumann's boundary conditions while a similar error bound could be obtained using Thomas' method and  $2^I$  points.

For cyclic reduction on Neumann's boundary conditions, if  $I \geq 6$  then an extra Fourier analysis and synthesis must be performed to solve the last column. This time should be divided by the number of rows and added to the time per row for cyclic reduction on Neumann's boundary conditions.

If odd/even reduction is used then the odd columns of  $U$  are calculated by solving tridiagonal systems. These systems are already restricted in size by the FFT performed on the even columns. One cyclic reduction program can be written to solve for  $U$  on the odd columns for all three boundary conditions. Thus cyclic reduction will be used to solve for  $U$  on the odd columns when odd/even reduction is employed.

### 3.4 Fourier Analysis and Synthesis

Hockney wrote one routine which performs Fourier analysis or synthesis for any of the boundary conditions [Hockney, Page 201]. His routine assumes that there is plenty of temporary storage available. Although this is true on most large serial machines ILLIAC IV has a relatively small amount of storage compared with its speed. Thus a method which is as fast as Hockney's Fourier routine but does not require extra storage would be

Preferred. Cooley, the author of FFT, presents such a method in [Cooley, Page 320-323]. Cooley shows that it is possible to use the complex fast Fourier transform to obtain the sine series (required for Dirichlet boundary conditions), cosine series (required for Neumann boundary conditions) and real series (required for periodic boundary conditions). Both Hockney's and Cooley's methods place restrictions on the interior mesh sizes:  $2^I - 1$  for Dirichlet,  $2^I$  for periodic and  $2^I + 1$  for Neumann.

Cooley's algorithm for Dirichlet boundary conditions goes as follows: let  $M = 2^I$ ,  $Y(j)$  for  $j = 1, \dots, M-1$  be the coefficients of the Fourier sine series,  $b(k) = \sum_{j=1}^{M-1} Y(j) \sin \frac{\pi j k}{M}$ .

$$Y(0) = Y(M) = 0, \text{ and } Y(j) = -Y(2M-j) = -Y(-j) \text{ for } j = 0, \dots, M.$$

Define

$$X(j) = -[Y(2j+1) - Y(2j-1)] + Y(2j)i \text{ for } j = 0, \dots, M/2.$$

Thus

$$X(0) = -[Y(1) + Y(-1)] + Y(0)i = -2Y(1) + 0i$$

and

$$X(M/2) = -[Y(M+1) - Y(M-1)] + Y(M)i = 2Y(M-1) + 0i.$$

For  $j = 1, \dots, M/2-1$ ,

$$\begin{aligned} X(M/2+j) &= -Y(M+2j+1) + Y(M+2j-1) + Y(M+2j)i = Y(M-2j-1) - Y(M-2j+1) \\ &\quad - Y(M-2j)i. \end{aligned}$$

Let  $\tilde{X}(j) = C(j)$  and calculate  $A_1(j)$  and  $A_2(j)$  for  $j = 0, 1, \dots, M/4$  using

$$A_1(j) = C(j) + C(M/2+j) \text{ and } A_2(j) = [C(j) - C(M/2+j)]W_M^j$$

where

$$W_M^j = \cos \frac{2\pi j}{M} + i \sin \frac{2\pi j}{M}.$$

Then calculate  $A(j)$  and  $A(M/4+j)$  for  $j = 0, 1, \dots, M/4-1$



where

$$A(j) = A_1(j) + i A_2(j), \quad \tilde{A}(M/4+j) = A_1(j) - i A_2(j).$$

Thus, for  $j = 0, 1, \dots, M/4$ ,

$$\begin{aligned} A(j) = & -Y(2j+1) + Y(2j-1) + Y(M-2j-1) - Y(M-2j+1) + \sin \frac{2\pi j}{M} [Y(2j+1) \\ & - Y(2j-1) + Y(M-2j-1) - Y(M-2j+1)] + \cos \frac{2\pi j}{M} [Y(2j) + Y(M-2j)] \\ & + i \left\{ Y(M-2j) - Y(2j) + \sin \frac{2\pi j}{M} [Y(2j) + Y(M-2j)] + \right. \\ & \left. \cos \frac{2\pi j}{M} [Y(2j-1) - Y(2j+1) - Y(M-2j-1) + Y(M-2j+1)] \right\} \end{aligned} \quad (48.1)$$

and for  $j = 1, \dots, M/4$

$$\begin{aligned} A(M/2-j) = & Y(2j-1) - Y(2j+1) - Y(M-2j+1) + Y(M-2j-1) - \\ & \sin \frac{2\pi j}{M} [Y(2j+1) - Y(2j-1) + Y(M-2j-1) - Y(M-2j+1)] - \\ & \cos \frac{2\pi j}{M} [Y(2j) + Y(M-2j)] - i \left\{ Y(M-2j) - Y(2j) - \right. \\ & \sin \frac{2\pi j}{M} [Y(2j) + Y(M-2j)] - \cos \frac{2\pi j}{M} [Y(2j-1) - \\ & \left. Y(2j+1) - Y(M-2j-1) + Y(M-2j+1)] \right\}. \end{aligned} \quad (48.2)$$

Now the complex fast Fourier transform is applied to  $A$  giving

$$X(j) = \sum_{k=0}^{N/2-1} A(k) W_N^{jk}, \quad \text{for } j = 0, 1, \dots, M/2-1.$$

For  $j = 1, 3, 5, \dots, M-3$  let

$$\begin{aligned} b(j) &= X\left(\frac{j-1}{2}\right) \text{ imaginary and} \\ b(j+1) &= X\left(\frac{j+1}{2}\right) \text{ real; } b(M-1) = X\left(\frac{M}{2}-1\right) \text{ imaginary.} \end{aligned}$$

Finally  $b(j)$  for  $j = 1, \dots, M-1$  is calculated

$$4b(j) = [b(j) - b(M-j)] - [b(j) + b(M-j)] / [2 \sin \frac{\pi j}{M}]. \quad (49)$$

$$\text{Now we have } 4b(j) = 4 \sum_{k=1}^{M-1} Y(k) \sin \frac{\pi j k}{M}.$$

Fourier analysis calculates  $\frac{2}{N} b(j)$  and Fourier synthesis calculates  $b(j)$ .

Since we are dealing with linear equations we can multiply  $\rho$  in (1) and



INITIAL CONDITIONS								
REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS	
\$X	PI	\$CO	I	\$DO	J	\$D3	KP	
\$R	P[\$X](2)	\$C2	I+ILJ	\$D1	L	\$D4	KM	
\$S	P[\$X](3)	\$C3	J-I+ILJ	\$D2	K	\$D5	ILJ	
LOOP	ASSIGNMENT STATEMENTS		OPERATIONS AND OVERLAP			TIME IN CLOCKS	TOTAL TIME FOR LOOP	
1	P[\$X](3):=\$R-\$S		PEM Store, Addition			14	$30\left(\frac{M-1}{2}\right)\left[\frac{N}{64}\right]$	
1	\$C3:=\$C3-\$D2		CU Addition overlapped			0		
1	\$B:=P[\$X](2)		PE Fetch overlapped			0		
1	\$A:=\$S;\$S:=\$B		Two PE Register Transfers			2		
1	P[\$X](2):=\$R+\$A		Addition, PE Store			14		
1	\$C2:=\$C2+\$D2		CU Addition overlapped			0		
1	\$R:=P[\$X](2)		PE Fetch overlapped			0		
-----								
	P[\$X](2):=\$R+\$R		Addition, PE Store			14	$42\left[\frac{N}{64}\right]$	
	\$C2:=\$DO+\$D5-\$D2		CU Addition overlapped			0		
	\$S:=P[\$X]-(2)		PE Fetch overlapped			0		
	\$C3:=\$D5		CU Register Transfer overlapped			0		
	P[\$X](3):,\$S-\$S		Addition, PE Store			14		
	\$C3:=\$C3+\$D2		CU Addition overlapped			0		
	\$R:=P[\$X](3) %ODD3		PE Fetch overlapped			0		
	P[\$X](3):=-\$R-\$R		Addition, PE Store			14		
	\$CO:=\$D3		CU register transfer overlapped			0		
-----								
2	\$C3:=\$C3+\$D3;\$C2:=\$C3-\$D2		CU Addition overlapped			0	$159\left(\frac{M-1}{4}-1\right)\left[\frac{N}{64}\right]$	
2	\$R:=P[\$X](3)-\$R		PE Fetch overlapped, Addition			7		
2	\$D6:=GRABONE(DS[GL],N2-IL)		CU Load			10		
2	\$D7:=GRABONE(DS[GL],IL)		CU Load			10		
2	\$S:=\$D6*\$R		Multiplication			9		
2	\$A:=P[\$X](2)*\$D7		PE Fetch overlapped, Multiplication					
2	\$S:=\$A-\$S %ODD2		Addition			7		
2	\$R:=\$R*\$D7		Multiplication			9		
2	\$A:=P[\$X](2)*\$D6		PE Fetch overlapped			9		
2	\$R:=\$A+\$R %ODD1		Addition			7		
2	\$C3:=\$DO+\$D4-\$CO		CU Addition overlapped			0		
2	\$CO:=\$CO+\$D3;\$C1:,\$C3-\$D3		CU Addition overlapped			0		
2	P[\$X](2):=P[\$X](3)-P[\$X](1)		Two PE Fetches, One overlapped					
			Addition, PE Store			21		
2	\$C2:=\$C2+\$D2;\$C1:=\$C1+\$D2		CU Addition overlapped			0		
2	ODD3:=P[\$X](2)		PE Fetch overlapped, PE Store			7		
2	P[\$X](2):=\$S-P[\$X](1)		PE Fetch, Addition, PE Store			21		
2	P[\$X](1):=\$S+P[\$X](1)		PE Fetch overlapped, Addition, PE Store			14		
2	\$C2:=\$C2-\$D2		CU Addition overlapped			0		
2	\$S:=P[\$X](2)		PE fetch overlapped			0		
2	P[\$X](1):=\$S-\$R		Addition, PE Store			14		
2	\$B:=\$R		PE Register Transfer overlapped			0		
2	\$R:=ODD3		PE Fetch overlapped			0		
2	P[\$X](2):=\$S+\$B		Addition, PE Store			14		
-----								
	\$C2:=\$D4+\$D1;\$C3:=\$D4+\$D5		CU Addition overlapped			0		$28\left[\frac{N}{64}\right]$
	\$R:=P[\$X](2)		PE Fetch overlapped			0		
	P[\$X](2):=P[\$X](3)		PE Fetch, PE Store			14		
	P[\$X](3):=\$R+\$R		Addition, PE Store			14		
SUBTOTAL						$(57M-155)\left[\frac{N}{64}\right]$		

Table 9. Preparation of the Data for the Fast Fourier Transform\*

\*The logic for the code to keep current values in the ADB's defined in the initial conditions is supplied in GLYPNIR, Appendix B.

INITIAL CONDITIONS									
REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS		
\$X	PI	\$D2	K	\$D5	ILJ	\$D8	I2		
\$DO	J	\$D3	KP	\$D6	IP	\$D9	CJ		
\$D1	L	\$D4	KM	\$D7	J1	\$D10	IT		
LOOP	ASSIGNMENT STATEMENTS		OPERATIONS AND OVERLAP		TIME IN CLOCKS	TOTAL TIME FOR LOOP			
	\$C2;=\$D5		CU Register transfer overlapped		0	0			
3	\$C3:=\$C2+\$D1		CU Addition overlapped		0	$35(\frac{M+1}{2})\lceil\frac{N}{64}\rceil$			
3	\$S:=P[\$X](2)		PE Fetch overlapped		0				
3	P[\$X](2):=\$S+P[\$X](3)		PE Fetch, Addition, PE Store		21				
3	P[\$X](3):=\$S-P[\$X](3)		PE Fetch overlapped, Addition, PE Store		14				
3	\$C2:=\$C2+\$D2		CU Addition overlapped		0				
	\$C0:=\$D3		CU Register Transfer overlapped		0	$10(\frac{M-5}{2})$			
	\$C2:=GRABONE(INDEX[GL],IL);		CU Load		10				
4	\$R:=P(0)		PE Fetch		7	$28(\frac{M-5}{2})\lceil\frac{N}{64}\rceil$			
4	P(0):=P(2)		PE Fetch, PE Store		14				
4	P(2):=\$R		PE Store		7				
4	\$C0;=\$C0+1;\$C2;=\$C2+1		CU Addition overlapped		0				
	\$D11:=GRABONE(DS[GL],N2-N11)		CU Load; Multiplication		19	$29(M-1)$			
	*\$D8				10				
	\$D12:=GRABONE(DS[GL],N11)		CU Load		10				
5	\$C2:=\$D4+\$D7+\$D5		CU Addition overlapped		0	$106(\text{Log}_2(M)-2)$ $(\frac{M+1}{8})\lceil\frac{N}{64}\rceil$			
5	\$C3:=\$C2+\$D9+\$D9;\$C1:=\$C3+\$C2		CU Addition overlapped		0				
5	\$R:=P[\$X](3)*\$D11		PE Fetch overlapped, Multiplication		9				
5	\$S:=P[\$X](1)		PE Fetch overlapped		0				
5	\$B:=\$S*\$D12		Multiplication		9				
5	\$R:=\$R-\$B %ODD1		Addition		7				
5	\$S:=\$S*\$D11		Multiplication		9				
5	\$A:=P[\$X](3)*\$D12		PE Fetch overlapped, Multiplication		9				
5	\$S:=\$A+\$S %ODD2		Addition		7				
5	P[\$X](3):=P[\$X](2)-\$R		PE Fetch overlapped, Addition, PE Store		14				
5	P[\$X](2):=P[\$X](2)+\$R		PE Fetch overlapped, Addition, PE Store		14				
5	\$C2:=\$C2+\$D2		CU Addition overlapped		0				
5	P[\$X](1):=P[\$X](2)-\$S		PE Fetch overlapped, Addition, PE Store		14				
5	P[\$X](2):=P[\$X](2)+\$S		PE Fetch overlapped, Addition, PE Store		14				
SUBTOTAL					$(\frac{53}{4}(M+1)\text{Log}_2(M)+5M-79)\lceil\frac{N}{64}\rceil+34M-54$				

Table 10. The Complex Fast Fourier Transform\*

\* The logic for the code to keep current values in the ADB's defined in the initial conditions is supplied in GLYPNIR, Appendix B.

INITIAL CONDITIONS							
REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS
\$X	PI	\$DO	J	\$D1	I	\$D2	ILJ
LOOP	ASSIGNMENT STATEMENT	OPERATIONS AND OVERLAP				TIME IN CLOCKS	TOTAL TIME FOR LOOP
	\$D3:=GRABONE(IS[GL],I1)	CU Load				10	$10(\frac{M+1}{2})$
-----							
6	\$C1:=\$D2+\$D1;\$C2:=\$D0-\$D1+\$D2	CU Addition overlapped				0	
6	\$R:=P[\$X](1)	PE Fetch overlapped				0	
6	\$B:=P[\$X](2)	PE Fetch				7	
6	\$R:=\$R=\$B	Addition				7	
6	\$A:=P[\$X](1)	PE Fetch overlapped				0	
6	\$C3:=\$D3	CU Register transfer overlapped				0	
6	\$A:=\$A+\$B	Addition				7	
6	\$S:=\$A*\$C3	Multiplication				9	
6	P[\$X](1):=\$R+\$S	Addition, PE Store				14	
6	P[\$X](2):=\$S-\$R	Addition, PE Store				14	$58(\frac{M+1}{2})\lceil\frac{N}{64}\rceil$
SUBTOTAL						$29M\lceil\frac{N}{64}\rceil+29\lceil\frac{N}{64}\rceil+5M+5$	

Table 11. Final Step in Cooley's Fourier Analysis and Synthesis\*

\*The logic for the code to keep current values in the ADB's defined in the initial conditions is supplied in GLYPNIR, Appendix B.

INITIAL CONDITIONS							
REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS	REGISTER	CONTENTS
\$X	PI	\$D1	IL	\$D3	L3	\$D5	L4
\$D0	CI	\$D2	L2	\$D4	K	\$D6	KM1
LOOP	ASSIGNMENT STATEMENTS		OPERATIONS AND OVERLAP		TIME IN CLOCKS		TOTAL TIME FOR LOOP
7	\$C2:=\$D1;\$C3:=\$D0		CU Register Transfer overlapped		0		20(M-1) $\left[\frac{N}{64}\right]$
7	P(2):=RTR(\$C3,,P(2));		PE Fetch, Route, PE Store		20		
-----			-----		-----		-----
	\$C1:=0;\$C3:+\$D0		CU Register clear overlapped		0		72 $\left[\frac{M}{64}\right]$
	\$S:=1.0/A(3)		PE Fetch, division		63		
	\$B:=P[\$X]*\$S		PE Fetch overlapped, Multiplication		9		
-----			-----		-----		-----
8	\$C1:=\$C1+1		CU Addition overlapped				94(N-1) $\left[\frac{M}{64}\right]$
8	\$A:=RTR(\$C1,,A(3))		PE Fetch, overlapped, Route		6		
8	\$S:=RTR(1,,\$S)		Route		3		
8	\$X:=RTR(1,,\$X)		Route		3		
8	\$R:=RTR(1,,\$B)		Route		3		
8	\$A:=\$A-\$S		Addition		7		
8	\$S:=1.0/\$A %E		Division		56		
8	\$A:=P[\$X]-\$R		PE Fetch overlapped, Addition		7		
8	\$B:=\$A*\$S %Q		Multiplication		9		
-----			-----		-----		
	\$A:=RTR(\$C1,,A(3))		PE Fetch overlapped, Route		6		22 $\left[\frac{M}{64}\right]$
	\$R:=\$B		PE Register transfer overlapped		0		
	\$A:=\$A*\$B		Multiplication		9		
	\$C2:=\$D6		CU Register transfer overlapped		0		
	\$S:=P[\$X](2)-\$A		PE Fetch overlapped, Addition		7		
	P[\$X](2):=\$R		PE Store overlapped		0		
-----			-----		-----		-----
9	\$C2:=\$D2+\$D3;\$C1:=\$C1-1		CU Addition overlapped		0		46(M-1) $\left[\frac{M}{64}\right]$
9	\$B:=RTL(1,,\$R) %Q		Route		3		
9	\$S:=RTL(1,,\$S) %E		Route		3		
9	\$X:=RTL(1,,\$X) %PI		Route		3		
9	\$A:=RTR(\$C1,,A(3))		PE Fetch, partially overlapped				
			Route		10		
9	\$R:=\$B %Q		PE Register transfer overlapped		0		
9	\$A:=\$A*\$S		Multiplication		9		
9	\$A:=P[\$X](2)-\$A		PE Fetch overlapped, Addition		7		
9	\$A:=\$A+\$R %T		Addition		7		
9	P[\$X](2):=\$S		PE Store overlapped		0		
9	\$B:=\$S %Q		PE Register transfer		1		
9	\$S:=\$A %E		PE Register transfer		1		
-----			-----		-----		
10	%C2:=\$D1;\$C3:=\$D0		CU Register transfer overlapped		0		20(M-1) $\left[\frac{N}{64}\right]$
10	P(2):=RTL(\$C3,,P(2))		PE Fetch, Route, PE Store		20		
-----			-----		-----		-----
SUBTOTAL					14 ON $\left[\frac{M}{64}\right]$	-46 $\left[\frac{M}{64}\right]$	+40(M-1) $\left[\frac{N}{64}\right]$

Table 12. CRED\*

\*The logic for the code to keep current values in the ADB's defined in the initial conditions is supplied in GLYPNIR, Appendix B.

the boundary conditions by  $\frac{1}{8M}$ . Then apply (48.1), (48.2), the complex fast Fourier transform and (49) before and after CRED. This algorithm, MFACR, is a modified FACR which solves (39) for Dirichlet's boundary conditions without using odd/even reduction.

### 3.5 Implementation of MFACR

We will discuss MFACR in explaining how it can be programmed in ASK and how much time such a routine would take. Appendix F contains GLYPNIR code of the algorithm. Appendix A contains information on timing methodology and explains the notation used in Tables 9 through 12. Both  $\Psi$  and  $U$  are  $M+2$  by  $N+2$  mesh.  $\Psi$  and  $U$  are stored in straight storage in  $C$  and  $P$ , respectively. MFACR consists of five steps:

1. Set up storage area  $P$
2. Fourier analysis uses  $\Psi$  to calculate  $\Psi^S$
3. CRED uses  $\Psi^S$  to calculate  $U^S$
4. Fourier synthesis uses  $U^S$  to calculate  $U$
5. Restore boundary conditions and  $C$ .

Step one places a linear combination of the boundary conditions and  $\Psi$  in the interior of the storage area  $P$ . In the process the first row of  $U$  along with the second and  $N+1$ st row of  $\Psi$  are lost. Thus they are saved in temporary storage. This section of the algorithm takes  $(23M+102)\left\lceil \frac{N}{64} \right\rceil + 13M$  clocks in ASK. Section 1 of Appendix F contains the GLYPNIR code.

Fourier synthesis and analysis are performed on up to 64 columns simultaneously described above. They are presented in three parts. First equations (48.1) and (48.2) prepare the data for the complex



fast Fourier transform. Table 9\* contains suggestions on how these equations could be programmed in ASK and supplies  $(57M-155) \left\lceil \frac{N}{64} \right\rceil$  clocks as a time estimate for this step. Table 10 contains suggestions on how the complex fast Fourier transform could be programmed in ASK. A time estimate of  $(\frac{53}{4}(M+1) \log_2(M+1) + 5M-79) \left\lceil \frac{N}{64} \right\rceil + 34M - 54$  is given for this step. Finally the implementation of equation (49) in ASK is discussed in Table 11 where a time estimate of  $29(M+1) \left\lceil \frac{N}{64} \right\rceil + 5M + 5$  is presented. The total time estimate for Cooley's algorithm is  $\{\frac{53}{4}(M+1) \log_2(M+1) + 91M - 205\} \left\lceil \frac{N}{64} \right\rceil + 39M - 49$  clocks. The GLYPNIR code for Cooley's method is found in section 2 of Appendix F.

MFACR uses the modified Thomas' method, equations (44), (45), and (47) to perform CRED on up to 64 rows simultaneously. Table 12 suggests how this method could be programmed in ASK and estimates how long that code would take,  $40(M-1) \left\lceil \frac{N}{64} \right\rceil + (140N-46) \left\lceil \frac{M}{64} \right\rceil$  clocks. Section 3 of Appendix F contains the GLYPNIR code for CRED.

The final step of MFACR consists of restoring the boundary conditions and the portions  $\Psi$  which have been altered. Section 4 of Appendix F contains the GLYPNIR code for this step. If programmed in ASK this step would take approximately  $42 \left\lceil \frac{N}{64} \right\rceil$  clocks.

The complete MFACR requires  $\{53/2(M-1) \log_2(M+1) + 245M-306\} \left\lceil \frac{N}{64} \right\rceil + (140N-46) \left\lceil \frac{M}{64} \right\rceil + 91M - 98$  clocks. The timing algorithm is taken from Appendix A.

---

\*

The loop numbers in Tables 9 to 12 correspond to the loop number in the GLYPNIR code in Appendix F. They are supplied to enable the reader to compare loops in the GLYPNIR code with the corresponding loops in the ASK code.

### 3.6 Implementation of FACR

FACR, one level of odd/even reduction would require  $(53/2(M-1) \log_2 (M+1) + 405M - 466) \left\lceil \frac{N}{128} \right\rceil + (80M + 104) \left\lceil \frac{N}{64} \right\rceil + (70N - 23) \left\lceil \frac{M}{64} \right\rceil + 81M - 84$  clocks. This figure assumes the odd columns are solved by Thomas' method with the e's calculated and stored in preconditioning. Thus the odd columns would take  $(84M - 46) \left\lceil \frac{N}{128} \right\rceil$  clocks. Setting up the even columns would take  $79M \left\lceil \frac{N}{128} \right\rceil$ . Preparing the odd rows for solution would take  $55M \left\lceil \frac{N}{128} \right\rceil$  clocks. CRED would take  $(70N-23) \left\lceil \frac{M}{64} \right\rceil$  clocks. Fourier analysis and synthesis would take  $(53/2(M-1) \log_2 (M+1) + 192M - 420) \left\lceil \frac{N}{128} \right\rceil + 68M - 48$ . The extra storage manipulation would take  $17M \left\lceil \frac{N}{64} \right\rceil$  clocks. The rest of the algorithm is identical to MFACR.

### 3.7 Summary of the Results on Direct Methods

In programming an algorithm on ILLIAC IV, an attempt is made to adapt the restrictions of the algorithm to maximize storage and machine efficiency. For efficient storage on ILLIAC IV, rows stored across PEM should contain a multiple of 64 words. To maximize machine efficiency the number of values being computed simultaneously equals 64. Hockney's method performs a Fourier analysis and synthesis on each column and performs CRED on each row. Thus to maximize machine efficiency on ILLIAC IV the number of interior rows and the number of interior columns must be multiples of 64. Fourier analysis and synthesis using Cooley's method requires  $2^I - 1$  points per column for Dirichlet's and Neumann's boundary conditions. If  $I \geq 6$  then storing the columns across PEM would waste 63 words per column. Thus we will store each column in one PE. Now we have columns restricted to  $2^I + 1$  ( $2^I$ ) for



Dirichlet's or Neumann's (periodic) boundary conditions. Rows are restricted to multiples of 64 to maximize storage efficiency. For periodic boundary conditions cyclic reduction saves time and temporary storage over Thomas' method [Hockney, Page 150], and restricts the mesh to  $2^I$  columns. This restriction on mesh size still allows maximum storage and machine efficiency. Cyclic reduction restricts the mesh to  $2^I + 1$  for Dirichlet's and Neumann's boundary conditions. This restriction reduces storage efficiency for both boundary conditions and it reduces machine efficiency for Neumann's boundary conditions. Thomas' method and the modified Thomas' method have no restrictions on the number of columns. Thomas' method is faster than the modified Thomas' method but requires temporary storage. Refer to Table 8 for a more detailed comparison of the methods available for use by CRED.

Hockney's FACR program handles the 9 different combinations of boundary conditions and performs one level of odd/even reduction in all cases. To maximize efficiency of such a program on ILLIAC IV different algorithms would be required for different mesh sizes and combinations of boundary conditions.

As stated earlier, odd/even reduction saves computer time on ILLIAC IV for certain mesh sizes but not for others. Thus FACR on ILLIAC IV should have two algorithms, one which uses odd/even reduction and one that does not. Furthermore, the choice of a method to solve CRED depends on the storage requirements of the individual problem. It is the author's opinion that modified Thomas' Method is the best compromise between storage and speed for Neumann's and Dirichlet's boundary conditions, and cyclic reduction is the best for periodic boundary conditions.

If Fourier analysis and synthesis is applied to columns which have Neumann's boundary conditions then there must be  $2^I + 1$  interior rows to be evaluated by CRED. This means CRED must be executed  $\left\lceil \frac{2^I + 1}{64} \right\rceil$  times. Note  $\left\lceil \frac{2^I + 1}{64} \right\rceil = \frac{2^I}{64} + 1$  if  $I \geq 6$  and thus for one execution of CRED, ILLIAC IV is working at  $1/64^{\text{th}}$  of its capacity. The user can avoid this machine inefficiency by setting up the program so that Fourier analysis and synthesis is performed in the direction which has non-Neumann boundary conditions.

On a M by N mesh the Fourier part of FACR requires the order of  $NM \log_2 M$  clocks when it is applied to the N columns in the mesh. CRED requires the order of NM clocks. By letting M be the smaller of the dimensions the user saves computer time.

#### 4. USE OF ILLIAC DISK FOR NON-CORE CONTAINED MESHES

When the problem being run becomes too large to contain in core, the I/O time becomes an important consideration. First the I/O system for ILLIAC IV will be discussed and then its efficiency for a specific problem will be examined.

##### 4.1 ILLIAC IV I/O System

The ILLIAC IV Disk is a 15,600 K word memory. This memory is divided into 52 bands. Each band is divided into 300 pages, each containing 16 lines of PEM (1 K words). Memory transfers between the ILLIAC IV Disk and PEM are restricted to pages. A data request can read or write up to 128 consecutive pages on one band. If the data is not consecutive or if it is spread over more than one band a separate request is needed for each string of consecutive pages. The time required for the disk to prepare to perform a data request is equivalent to the time to transfer two pages of data between the Disk and PEM. Thus if a data request reads or writes on page  $i$  and band  $j$ , the next data request should skip at least two pages and start at page  $i + 3$  of band  $k$ , where  $j$  need not equal  $k$ . If the second data request wanted page  $i + 1$ , or  $i + 2$ , the request would have to wait for a revolution of the disk before it could be executed. The transfer rate between PEM and Disk is 133 usec per page and the disk revolves once every 40 msec. One revolution of the disk is equivalent to 640,000 ILLIAC IV clocks.

##### 4.2 I/O for the Bernard-Rayleigh convection problem

Now let us examine the Bernard-Rayleigh convection problem where the following equations are used to solve for temperature,  $T$ , pressure,  $P$ , and velocity components  $u$  and  $w$  along with the  $X$  and  $Z$  axes:

$$\frac{\partial T}{\partial t} + u \frac{\partial T}{\partial x} + w \frac{\partial T}{\partial z} = \frac{1}{P_r} \nabla^2 T \quad (50)$$

$$\frac{\partial \eta}{\partial t} + u \frac{\partial \eta}{\partial x} + w \frac{\partial \eta}{\partial z} = \frac{R_a}{P_r} \frac{\partial T}{\partial x} \nabla^2 \eta. \quad (51)$$

$$\eta = \frac{\partial w}{\partial x} - \frac{\partial u}{\partial z} = \nabla^2 \psi \quad (52)$$

where  $t$  is time,  $R_a$  is the Rayleigh number,  $P_r$  is the Prandtl number, and  $\psi$  is the stream function defined by

$$u = - \frac{\partial \psi}{\partial z}, \quad w = \frac{\partial \psi}{\partial x}$$

The finite difference schemes employed to solve these equations require five meshes:  $T^{(\tau)}$ ,  $T^{(\tau-1)}$ ,  $\eta^{(\tau)}$ ,  $\eta^{(\tau-1)}$  and  $\psi^{(\tau)}$ .

Each time step of the convection process has three main parts;  
calculate  $\eta^{(\tau+1)}$ , calculate  $T^{(\tau+1)}$  and calculate  $\psi^{(\tau+1)}$ .

$\eta_{i,j}^{(\tau+1)}$  is calculated using  $\eta_{i+1,j}^{(\tau)}$ ,  $\eta_{i,j+1}^{(\tau)}$ ,  $\eta_{i,j}^{(\tau-1)}$ ,  $T_{i+1,j}^{(\tau)}$ , and  $\psi_{i+1,j+1}^{(\tau)}$ .  $T_{i,j}^{(\tau+1)}$

is calculated using  $T_{i+1,j}^{(\tau)}$ ,  $T_{i,j+1}^{(\tau)}$ ,  $T_{i,j}^{(\tau-1)}$ ,  $\psi_{i+1,j+1}^{(\tau)}$ ,  $\psi_{i+1,j}^{(\tau)}$  and  $\psi_{i,j+1}^{(\tau)}$ .

Using all of  $\eta^{(\tau+1)}$ ,  $\psi_{i,j}^{(\tau+1)}$  is calculated.

The process will be divided into two parts. First the prognostic equations (50) and (51) are solved to obtain  $\eta^{(\tau+1)}$  and  $T^{(\tau+1)}$ . Then Poisson's equation, (52), is solved for  $\psi^{(\tau+1)}$ .

The prognostic equations require five meshes while Poisson's equation requires two meshes if solved by an iterative method and one mesh if solved by a direct method. Thus non-core contained problems can be divided into two types. In the first type, the meshes required for solving Poisson's equation are core contained but some of the meshes required for solving the prognostic equations are not contained in core.

In the second case, the meshes are so large that none of the meshes can be core contained.

In [Ogura, et al] the first type of problem was studied in which SOR was used to solve Poisson's equation. If a direct method had been used in that study, the maximum mesh size could be doubled.

In this study the second type of I/O problem will be examined where MFACR is used to solve Poisson's equation. For an example we will use 512 by 512 meshes.  $P_{i,j}$  where  $1 \leq i \leq 64$  and  $0 \leq j \leq 3$  is a page of a mesh containing all points  $a_{k,\ell}$  where  $8_i - 8 \leq k < 8_i$  and  $128 j \leq \ell < 128 j + 128$ .

If the prognostic equations are solved separately using ASK code they are found to be about 50% I/O bound. To decrease the I/O bound the Fourier analysis of MFACR will be performed along with the calculation of the prognostic equations.

This will be followed by CRED, with the final step being Fourier synthesis.

The data will be stored in blocks 22 pages long. Block number I,  $B_I$  contains  $P_{I,j}$  where  $0 \leq j \leq 3$  of  $\eta^{(\tau)}$ ,  $T^{(\tau)}$ ,  $\eta^{(\tau-1)}$ ,  $T^{(\tau-1)}$ , and  $\psi^{(\tau)}$  for time step  $\tau$ . Also included in  $B_I$  are the two pages required for the read or write interpret. Thirteen blocks can be placed on one band with 14 pages remaining. The data will be spread across the 11 bands of disk with 22 pages separating each block to leave room for the information from PEM to be written. Seven blocks of data and six blocks for writing are found on bands 0, 2, 4, 6, and 8. Six blocks of data and seven blocks for writing are found on bands 1, 3, 5, and 7. Band 9 contains 5 blocks of data and seven blocks for writing. Band 10 contains the last block for writing.

The blocks are skewed to enable efficient access of rows and columns. The first block on band I starts at page  $-2I \bmod 300$ .

The prognostic equations and the Fourier analysis are performed while the disk revolves ten times and 110 pages are passed on the eleventh revolution. Figure 6 shows what data will be printed on the first 44 pages of each band.

Only one column of pages can fit in PEM at one time. Since CRED requires an entire column before the calculations can be performed, a column is read into PEM, the calculations are performed and the column is printed on disk. The data has been skewed in such a way as to enable an entire column to be read or written in one revolution, see Figure 6. There are four columns and the calculation of each takes  $1/4$  of a revolution. The total I/O time for CRED is nine revolutions.

The final step, Fourier Synthesis, is performed in four revolutions. Figure 6 shows that this can be accomplished on the first 44 pages by calculating rows 5, 31 and 57 on the first revolution, rows 18 and 44 on the second revolution, rows 12, 38, 64 on the third revolution and rows 25, 51 on the last revolution. To enable  $\Psi^{(\tau+1)}$  to overwrite  $\Psi^{(\tau+1/3)}$  Row I of  $\Psi^{(\tau+2/3)}$  is written 3 data blocks before Row I of  $\Psi^{(\tau+1/3)}$  during CRED.\*

The entire time step required 23 revolutions and 110 pages on the 24th revolution. This takes 932 milliseconds. The calculation could be performed in 333 milliseconds. Thus the process is 64% I/O bound.

A large portion of this I/O bound is created by the solution of Poisson's equation. The last two steps of MFACR required 13 revolutions of the disk while the calculations for those steps required the amount of time for  $2\frac{1}{2}$  revolutions.

\*  $\Psi^{(\tau+1/3)}$  contains the harmonics calculated by Fourier analysis.  
 $\Psi^{(\tau+2/3)}$  contains the harmonics calculated by CRED.



To minimize total computer time a problem takes, a programmer must adapt algorithms to disk maps which give minimum I/O time. With I/O being such a dominant factor in the above problem the use of library subroutines becomes minimal because programs need to be customized to minimize I/O.



Figure 6. A Disk Map of 512 by 512 Meshes

Page	Band 0	Band 1	Band 2	Band 3	Band 4	Band 5	Band 6	Band 7	Band 8	Band 9	Band 10
0		$T_{5,0}^{(\tau+1)}$		$\eta_{18,0}^{(\tau+1)}$		$T_{31,0}^{(\tau)}$		$\eta_{44,0}^{(\tau)}$		$\psi_{57,0}^{(\tau+\frac{1}{3})}$	
1		$T_{5,1}^{(\tau+1)}$		$\eta_{18,1}^{(\tau+1)}$		$T_{31,1}^{(\tau)}$		$\eta_{44,1}^{(\tau)}$		$\psi_{57,1}^{(\tau+\frac{1}{3})}$	
2		$T_{5,2}^{(\tau+1)}$		$\eta_{18,2}^{(\tau+1)}$		$T_{31,2}^{(\tau)}$		$\eta_{44,2}^{(\tau)}$		$\psi_{57,2}^{(\tau+\frac{1}{3})}$	
3		$T_{5,3}^{(\tau+1)}$		$\eta_{18,3}^{(\tau+1)}$		$T_{31,3}^{(\tau)}$		$\eta_{44,3}^{(\tau)}$		$\psi_{57,3}^{(\tau+\frac{1}{3})}$	
4		$\eta_{5,0}^{(\tau+1)}$		$T_{18,0}^{(\tau)}$		$\eta_{31,0}^{(\tau)}$		$\psi_{44,0}^{(\tau+\frac{1}{3})}$			$T_{64,0}^{(\tau+1)}$
5		$\eta_{5,1}^{(\tau+1)}$		$T_{18,1}^{(\tau)}$		$\eta_{31,1}^{(\tau)}$		$\psi_{44,1}^{(\tau+\frac{1}{3})}$			$T_{64,1}^{(\tau+1)}$
6		$\eta_{5,2}^{(\tau+1)}$		$T_{18,2}^{(\tau)}$		$\eta_{31,2}^{(\tau)}$		$\psi_{44,2}^{(\tau+\frac{1}{3})}$			$T_{64,2}^{(\tau+1)}$
7		$\eta_{5,3}^{(\tau+1)}$		$T_{18,3}^{(\tau)}$		$\eta_{31,3}^{(\tau)}$		$\psi_{44,3}^{(\tau+\frac{1}{3})}$			$T_{64,3}^{(\tau+1)}$
8		$T_{5,0}^{(\tau)}$		$\eta_{18,0}^{(\tau)}$		$\psi_{31,0}^{(\tau+\frac{1}{3})}$			$T_{51,0}^{(\tau+1)}$		$\eta_{64,0}^{(\tau+1)}$
9		$T_{5,1}^{(\tau)}$		$\eta_{18,1}^{(\tau)}$		$\psi_{31,1}^{(\tau+\frac{1}{3})}$			$T_{51,1}^{(\tau+1)}$		$\eta_{64,1}^{(\tau+1)}$
10		$T_{5,2}^{(\tau)}$		$\eta_{18,2}^{(\tau)}$		$\psi_{31,2}^{(\tau+\frac{1}{3})}$			$T_{51,2}^{(\tau+1)}$		$\eta_{64,2}^{(\tau+1)}$
11		$T_{5,3}^{(\tau)}$		$\eta_{18,3}^{(\tau)}$		$\psi_{31,3}^{(\tau+\frac{1}{3})}$			$T_{51,3}^{(\tau+1)}$		$\eta_{64,3}^{(\tau+1)}$
12		$\eta_{5,0}^{(\tau)}$		$\psi_{18,0}^{(\tau+\frac{1}{3})}$			$T_{38,0}^{(\tau+1)}$		$\eta_{51,0}^{(\tau+1)}$		$T_{64,0}^{(\tau)}$
13		$\eta_{5,1}^{(\tau)}$		$\psi_{18,1}^{(\tau+\frac{1}{3})}$			$T_{38,1}^{(\tau+1)}$		$\eta_{51,1}^{(\tau+1)}$		$T_{64,1}^{(\tau)}$
14		$\eta_{5,2}^{(\tau)}$		$\psi_{18,2}^{(\tau+\frac{1}{3})}$			$T_{38,2}^{(\tau+1)}$		$\eta_{51,2}^{(\tau+1)}$		$T_{64,2}^{(\tau)}$
15		$\eta_{5,3}^{(\tau)}$		$\psi_{18,3}^{(\tau+\frac{1}{3})}$			$T_{38,3}^{(\tau+1)}$		$\eta_{51,3}^{(\tau+1)}$		$T_{64,3}^{(\tau)}$
16		$\psi_{5,0}^{(\tau+\frac{1}{3})}$			$T_{25,0}^{(\tau+1)}$		$\eta_{38,0}^{(\tau+1)}$		$T_{51,0}^{(\tau)}$		$\eta_{64,0}^{(\tau)}$
17		$\psi_{5,1}^{(\tau+\frac{1}{3})}$			$T_{25,1}^{(\tau+1)}$		$\eta_{38,1}^{(\tau+1)}$		$T_{51,1}^{(\tau)}$		$\eta_{64,1}^{(\tau)}$
18		$\psi_{5,2}^{(\tau+\frac{1}{3})}$			$T_{25,2}^{(\tau+1)}$		$\eta_{38,2}^{(\tau+1)}$		$T_{51,2}^{(\tau)}$		$\eta_{64,2}^{(\tau)}$
19		$\psi_{5,3}^{(\tau+\frac{1}{3})}$			$T_{25,3}^{(\tau+1)}$		$\eta_{38,3}^{(\tau+1)}$		$T_{51,3}^{(\tau)}$		$\eta_{64,3}^{(\tau)}$
20			$T_{12,0}^{(\tau+1)}$		$\eta_{25,0}^{(\tau+1)}$		$T_{38,0}^{(\tau)}$		$\eta_{51,0}^{(\tau)}$		$\psi_{64,0}^{(\tau+\frac{1}{3})}$

21	$T_{12,1}^{(\tau+1)}$	$n_{25,1}^{(\tau+1)}$	$T_{38,1}^{(\tau)}$	$n_{51,1}^{(\tau)}$	$\psi_{64,1}^{(\tau+\frac{1}{3})}$
22	$T_{12,2}^{(\tau+1)}$	$n_{25,2}^{(\tau+1)}$	$T_{38,2}^{(\tau)}$	$n_{51,2}^{(\tau)}$	$\psi_{64,2}^{(\tau+\frac{1}{3})}$
23	$T_{12,3}^{(\tau+1)}$	$n_{25,3}^{(\tau+1)}$	$T_{38,3}^{(\tau)}$	$n_{51,3}^{(\tau)}$	$\psi_{64,3}^{(\tau+\frac{1}{3})}$
24	$n_{12,0}^{(\tau+1)}$	$T_{25,0}^{(\tau)}$	$n_{38,0}^{(\tau)}$	$\psi_{51,0}^{(\tau+\frac{1}{3})}$	
25	$n_{12,1}^{(\tau+1)}$	$T_{25,1}^{(\tau)}$	$n_{38,1}^{(\tau)}$	$\psi_{51,1}^{(\tau+\frac{1}{3})}$	
26	$n_{12,2}^{(\tau+1)}$	$T_{25,2}^{(\tau)}$	$n_{38,2}^{(\tau)}$	$\psi_{51,2}^{(\tau+\frac{1}{3})}$	
27	$n_{12,3}^{(\tau+1)}$	$T_{25,3}^{(\tau)}$	$n_{38,3}^{(\tau)}$	$\psi_{51,3}^{(\tau+\frac{1}{3})}$	
28	$T_{12,0}^{(\tau)}$	$n_{25,0}^{(\tau)}$	$\psi_{38,0}^{(\tau+\frac{1}{3})}$		$T_{58,0}^{(\tau+1)}$
29	$T_{12,1}^{(\tau)}$	$n_{25,1}^{(\tau)}$	$\psi_{38,1}^{(\tau+\frac{1}{3})}$		$T_{58,1}^{(\tau+1)}$
30	$T_{12,2}^{(\tau)}$	$n_{25,2}^{(\tau)}$	$\psi_{38,2}^{(\tau+\frac{1}{3})}$		$T_{58,2}^{(\tau+1)}$
31	$T_{12,3}^{(\tau)}$	$n_{25,3}^{(\tau)}$	$\psi_{38,3}^{(\tau+\frac{1}{3})}$		$T_{58,3}^{(\tau+1)}$
32	$n_{12,0}^{(\tau)}$	$\psi_{25,0}^{(\tau+\frac{1}{3})}$		$T_{45,0}^{(\tau+1)}$	$n_{58,0}^{(\tau+1)}$
33	$n_{12,1}^{(\tau)}$	$\psi_{25,1}^{(\tau+\frac{1}{3})}$		$T_{45,1}^{(\tau+1)}$	$n_{58,1}^{(\tau+1)}$
34	$n_{12,2}^{(\tau)}$	$\psi_{25,2}^{(\tau+\frac{1}{3})}$		$T_{45,2}^{(\tau+1)}$	$n_{58,2}^{(\tau+1)}$
35	$n_{12,3}^{(\tau)}$	$\psi_{25,3}^{(\tau+\frac{1}{3})}$		$T_{45,3}^{(\tau+1)}$	$n_{58,3}^{(\tau+1)}$
36	$\psi_{12,0}^{(\tau+\frac{1}{3})}$		$T_{32,0}^{(\tau+1)}$	$n_{45,0}^{(\tau+1)}$	$T_{58,0}^{(\tau)}$
37	$\psi_{12,1}^{(\tau+\frac{1}{3})}$		$T_{32,1}^{(\tau+1)}$	$n_{45,1}^{(\tau+1)}$	$T_{58,1}^{(\tau)}$
38	$\psi_{12,2}^{(\tau+\frac{1}{3})}$		$T_{32,2}^{(\tau+1)}$	$n_{45,2}^{(\tau+1)}$	$T_{58,2}^{(\tau)}$
39	$\psi_{12,3}^{(\tau+\frac{1}{3})}$		$T_{32,3}^{(\tau+1)}$	$n_{45,3}^{(\tau+1)}$	$T_{58,3}^{(\tau)}$
40		$T_{19,0}^{(\tau+1)}$	$n_{32,0}^{(\tau+1)}$	$T_{45,0}^{(\tau)}$	$n_{58,0}^{(\tau)}$
41		$T_{19,1}^{(\tau+1)}$	$n_{32,1}^{(\tau+1)}$	$T_{45,1}^{(\tau)}$	$n_{58,1}^{(\tau)}$
42		$T_{19,2}^{(\tau+1)}$	$n_{32,2}^{(\tau+1)}$	$T_{45,2}^{(\tau)}$	$n_{58,2}^{(\tau)}$
43		$T_{19,3}^{(\tau+1)}$	$n_{32,3}^{(\tau+1)}$	$T_{45,3}^{(\tau)}$	$n_{58,3}^{(\tau)}$



## 5. CONCLUSIONS

Table 13 compares MFACR and FACR with SOR and ADI for Dirichlet's boundary conditions. The direct methods use modified Thomas' method to perform CRED and Cooley's methods to perform the Fourier part of the algorithm. Unless an excellent initial guess is supplied for the iterative methods, the direct methods are much faster.

For maximum machine efficiency, MFACR requires  $64k$  by  $2^{I+1}$  meshes, FACR requires  $128k$  by  $2^{I+1}$  meshes, ADI requires  $64k$  by  $64I$  meshes, and SOR requires  $64k$  by  $j$  meshes. As seen in section 2, SOR can partition its mesh to obtain maximum machine efficiency. If the mesh for one of the other methods did not meet the row restrictions, the mesh could not be partitioned to meet the restrictions. A possible solution would be to solve a number of meshes at the same time.

MESH SIZE M+2 by N+2	FACR	MFACR	SOR PER ITERATION	ADI PER ITERATION
33 by 63	6.51 or 25300	5.71 or 22400	I or 3900	3.81 or 14900
65 by 63	5.31 or 49500	4.91 or 39100	I or 7900	2.51 or 20100
65 by 127	3.71 or 59500	4.71 or 74000	I or 15900	2.61 40700
129 by 127	3.81 or 122000	4.91 or 156000	I or 32000	2.71 or 87700

Table 13. A Comparison of Methods with Respect to Time\*

---

\*

Two numbers are supplied. The number times I is the number of iterations which could be performed in an equivalent amount of time. The other number is the number of clocks required on ILLIAC IV. The times given for SOR in section 2.3 don't take into account the time required to check for convergence. To calculate the times for SOR in Table 13, 20% has been added to the times in section 2.3 to account for convergence checking.

Direct methods require only one mesh in core. The mesh initially contains the interior source points and the boundary conditions. Throughout the algorithm, the mesh is used as temporary storage with the values in the mesh set equal to the solution in the final step. Iterative methods require both the source and the solution mesh at every iteration.

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APPENDIX A

## TIMING METHODOLOGY

The procedure for estimating the amount of time required by an algorithm is explained below. Care has been taken to insure that the method of calculating the computer time does not give one algorithm an unwarranted advantage over another.

In determining the amount of time an algorithm takes on a computer we must consider the language in which the algorithm is programmed. In comparing GLYPNIR to ASK, one finds that the average GLYPNIR code takes about twice as long as the ASK code for the same problem. Two major reasons for this difference are memory fetching and indexing. The code generated by GLYPNIR for indexing can be improved considerably by rewriting it in ASK. GLYPNIR does fetching and storing of data between FEM and PE registers that can be eliminated by efficient use of the PE registers in ASK.

GLYPNIR code is much easier to write than ASK. One must choose between simplicity and speed. However, GLYPNIR permits the programmer to write sections of code in ASK. Thus the programmer can write code in GLYPNIR and then rewrite the statements which are executed most often in ASK. In estimating the time for the different methods, we have timed the statements which are executed in the inner loops of the iterative method. Since these statements take most of the computer time they should be written in ASK. We use the GLYPNIR statements as an outline of the method but the time estimates reflect efficient ASK code, not the code that GLYPNIR would generate. To show the reader how we arrived at our time estimates we have included a table with assignment statements, operations which are not overlapped and the clocks required for each method.

The assignment statements are written in a combination of GLYPNIR and ASK. The parts of the statements which are not used in GLYPNIR are defined in Table 14. A summary of the assumptions used in obtaining the time estimates is found in Table 15.

NOTATION	MEANING	NOTATION	MEANING
\$A	PE Register A	\$Ci	ACAR i for $i=0,1,2,3$
\$B	PE Register B	U(i)	U indexed by ACAR i
\$R	PE Register R	U[\$X]	U indexed by register X
\$S	PE Register S	U[\$X](i)	both \$X and \$Ci indexing of U
\$X	PE Register X	\$Di	ADB storage location i for $i=0,\dots,63$

Table 14. Assignment Statement Notation

Every iterative algorithm we consider in this study will have two parts: Phase 1 will be used to improve 64 points simultaneously; Phase 2 will improve up to 63 points simultaneously. The machine efficiency of Phase 1 is greater than the machine efficiency of Phase 2. If one used only Phase 1 on certain meshes the extra time needed to bring the data to the PE's would exceed the time saved by improving machine efficiency. Thus the fastest code is obtained by a combination of Phase 1 and Phase 2.

- 1) CU instructions are overlapped completely by a combination of FINST/PE instructions and PEM fetches.
- 2) Memory fetching is overlapped as much as possible.
- 3) Seven PE clocks are used to load ((store) a PE register from (to) PEM.
- 4) One PE clock issued to load a PE register from the CU.
- 5) Transfer between PE registers require one clock.
- 6) Each GRABONE will be considered to be an ASK LOAD instruction (10 clocks).

Table 15. Assumptions in Timing\*

\*

The author feels these timing assumptions should give times within 20% of the actual times the algorithms would take on ILLIAC IV.

To aid in comparing times of different methods we will define an execution of Phase 1 to be the improvement of  $64$  points simultaneously by Phase 1. Similarly, an execution of Phase 2 is the improvement of  $k$  points where Phase 2 improves  $k$  points simultaneously.



APPENDIX B

MAJOR STEPS OF FACR AND MFACR FOR DIRICHLET'S BOUNDARY CONDITIONS

Given the meshes  $\Psi$  and  $U$  with mesh points at locations  $\Psi_{i,j}$  and  $U_{i,j}$  where  $0 \leq i \leq M$  and  $0 \leq j \leq N$  the five point difference scheme approximations of Poisson's equation becomes

$$\frac{U_{s-1,t} - 2U_{s,t} + U_{s+1,t}}{h_x^2} + \frac{U_{s,t-1} - 2U_{s,t} + U_{s,t+1}}{h_y^2} = \Psi_{s,t} \quad (B.1)$$

To solve (B.1) FACR works as follows.

Odd/even reduction (see Section 3.1) calculates  $\Psi^*$  for the even columns of  $\Psi$ . Equation (41) from Section 3.1 can be expanded into the following equation for individual mesh points.

$$\begin{aligned} & \frac{1}{h_y^2} (U_{i,t-2} + U_{i,t+2}) - \frac{h_x^2}{h_y^4} (U_{i-2,t} + U_{i+2,t}) \\ & + 4 \left( \frac{h_y^2}{h_x^4} + \frac{1}{h_x^2} \right) (U_{i-1,t} + U_{i+1,t}) - \left( 6 \frac{h_y^2}{h_x^4} + \frac{8}{h_x^2} + \frac{2}{h_y^2} \right) U_{i,t} = \Psi_{i,t}^* \end{aligned} \quad (B.2)$$

where

$$\Psi_{i,t}^* = \Psi_{i,t-1} - \frac{h_y^2}{h_x^2} (\Psi_{i-1,t} + \Psi_{i+1,t}) + 2 \left( \frac{h_y^2}{h_x^2} + 1 \right) \Psi_{i,t} + \Psi_{i,t-1} \quad (B.2.1)$$

Fourier analysis calculates the Fourier harmonics  $\Psi^S$  and  $\Psi^*$  defined by

$$\Psi_{i,t}^S = \frac{2}{M} \sum_{k=1}^{M-1} \Psi_{k,t}^* \sin \frac{\pi k i}{M}. \quad (B.3)$$

Combining (B.2) and (B.3) we obtain

$$\begin{aligned} \Psi_{i,t}^S = \frac{2}{M} \sum_{k=1}^{M-1} \left[ \sin \frac{\pi k i}{M} \left\{ \frac{1}{h_y^2} (U_{k,t-2} + U_{k,t+2}) - \frac{h_x^2}{h_y^4} (U_{k-2,t} + U_{k+2,t}) \right. \right. \\ \left. \left. + 4 \left( \frac{h_y^2}{h_x^4} + \frac{1}{h_x^2} \right) (U_{k-1,t} + U_{k+1,t}) - \left( 6 \frac{h_y^2}{h_x^4} + \frac{8}{h_x^2} + \frac{2}{h_x^2} \right) U_{k,t} \right\} \right] \end{aligned} \quad (B.4)$$

The Fourier harmonics for  $U$ ,  $U^S$  are defined as follows:

$$U_{i,t}^S = \frac{2}{M} \sum_{k=1}^{M-1} \text{SIN} \frac{\pi k i}{M} U_{k,t} \quad (\text{B.5})$$

and thus

$$U_{k,t} = \sum_{\ell=1}^{M-1} \text{SIN} \frac{\pi \ell k}{M} U_{\ell,t}^S. \quad (\text{B.6})$$

Combining (B.5), (B.6) and (B.4) we obtain

$$\begin{aligned} \Psi_{i,t}^S = & \frac{1}{h_y^2} (U_{i,t+2}^S + U_{i,t-2}^S) - \left( 6 \frac{h_y^2}{h_x^4} + \frac{8}{h_x^2} + \frac{2}{h_y^2} \right) U_{i,t}^S \\ & + \frac{2}{M} \sum_{k=1}^{M-1} \left[ \text{SIN} \frac{\pi k i}{M} \left\{ -\frac{h_y^2}{h_x^4} \sum_{\ell=1}^{M-1} \left( \text{SIN} \frac{\pi(k-2)\ell}{M} U_{\ell,t}^S + \text{SIN} \frac{\pi(k+2)\ell}{M} U_{\ell,t}^S \right) \right. \right. \\ & \left. \left. + 4 \left( \frac{h_y^2}{h_x^4} + \frac{1}{h_y^2} \right) \sum_{\ell=1}^M \left( \text{SIN} \frac{\pi(k-1)\ell}{M} U_{\ell,t}^S + \text{SIN} \frac{\pi(k+1)\ell}{M} U_{\ell,t}^S \right) \right\} \right]. \end{aligned} \quad (\text{B.7})$$

Using the facts that  $\text{Sin}(a \pm b) = \text{Sin } a \text{ Cos } b \pm \text{Sin } b \text{ Cos } a$  and

$$\sum_{k=1}^{M-1} \text{Sin} \frac{\pi \ell k}{M} \text{Sin} \frac{\pi i k}{M} = \frac{N}{2} \delta_{i,\ell} \text{ for } \ell, i = 1, 2, \dots, M-1 \text{ where}$$

$\delta_{i,\ell} = \begin{cases} 0 & \text{if } i \neq \ell \\ 1 & \text{if } i = \ell \end{cases}$  we can decouple (B.7) into a tridiagonal linear system

$$\begin{aligned} \Psi_{i,t}^S = & \frac{1}{h_y^2} (U_{i,t+2}^S + U_{i,t-2}^S) - \left( 2 \frac{h_y^2}{h_x^4} \text{Cos} \frac{2\pi i}{M} - 8 \left( \frac{h_y^2}{h_x^4} + \frac{1}{h_x^2} \right) \text{Cos} \frac{\pi i}{M} \right. \\ & \left. + 6 \frac{h_y^2}{h_x^4} + \frac{8}{h_x^2} + \frac{2}{h_y^2} \right) U_{i,t}^S \end{aligned} \quad (\text{B.8})$$

CRED solves (B.8) for  $U^S$ .

Fourier synthesis calculates  $U$  from  $U^S$  using (B.6).

This gives us the values for  $U$  on the even columns.

Using the values of  $U$  on the even columns we can solve (B.1) for the values of  $U$  on the odd columns using a method which solves tridiagonal linear systems.

In summary FACR performs the following algorithm:

1. Given  $\Psi$  compute  $\Psi^*$  for the even columns using (B.2.1) and overwrite  $\Psi$  with  $\Psi^*$  on the even columns.
2. Using  $\Psi^*$  compute  $\Psi^S$  according to (B.3) and overwrite  $\Psi^*$  with  $\Psi^S$ .
3. Using  $\Psi^S$  solve (B.8) for  $U^S$  and overwrite  $\Psi^S$  with  $U^S$ .
4. Using  $U^S$  compute  $U$  according to (B.6) on the even columns and overwrite  $U^S$  with  $U$  on the even columns.
5. Using the values of  $U$  on the even columns solve (B.1) for the values of  $U$  on the odd columns and overwrite  $\Psi$  with  $U$  on the odd columns.

MFACR is similar to FACR but the odd/even reduction step is omitted.

MFACR begins with Fourier analysis which calculates the Fourier harmonics  $\Psi^S$  for  $\Psi$  using

$$\Psi_{i,t}^S = \frac{2}{M} \sum_{k=1}^{M-1} \Psi_{k,t} \text{ SIN } \frac{\pi k i}{M} \quad (\text{B.9})$$

From B.9 we obtain

$$\Psi_{i,t}^S = \frac{2}{M} \sum_{k=1}^{M-1} \text{ SIN } \frac{\pi k i}{M} \left( \frac{U_{s-1,t} - 2U_{s,t} + U_{s+1,t}}{h_x^2} + \frac{U_{s,t-1} - 2U_{s,t} + U_{s,t+1}}{h_y^2} \right) \quad (\text{B.10})$$



Using the same methods employed to obtain (B.8) from (B.7) we obtain (B.11) from (B.10).

$$\psi_{i,t}^s = \frac{1}{h_y^2} (U_{i,t+1}^s + U_{i,t-1}^s) + \left( \frac{2}{h_y^2} \cos \frac{\pi i}{M} - \frac{2}{h_y^2} - \frac{2}{h_x^2} \right) U_{i,t}^s \quad (\text{B.11})$$

CRED solves (B.11) for  $U_{i,t}^s$ .

Finally (B.1) is used to calculate  $U$  for the entire mesh.

## SUCCESSIVE (POINT\_ OVER-RELAXATION METHOD IN GLYPNIR, MSOR

ILLIAC IV GLYPNIR COMPILER  
( 86500 VERSION 3.3.00 )

```

1  SUBROUTINE MSOR(PCPOINT OUT X,CREAL W8,CINT CM,CINT CM,CREAL BOUND.
2  PCPOINT EFA,CREAL DX,CREAL DZ,CINT OUT IT):
3
4  ***** DEFINITION OF PARAMETERS *****
5
6  X IS THE MESH TO BE IMPROVED
7
8  W8 IS THE OPTIMAL RELAXATION PARAMETER FOR SOR
9
10 W8 IS SUPPLIED BY THE MAIN PROGRAM
11
12 CM IS THE NUMBER OF ROWS IN THE MESH X
13
14 CM IS THE NUMBER OF COLUMNS IN THE MESH X
15
16 BOUND IS THE MAX ERROR AT WHICH ANOTHER ITERATION WILL
17 NOT BE PERFORMED
18
19 EFA IS THE CONSTANT MESH IN THE EQUATION AX=ETA OBTAINED
20 FROM THE POISSON'S EQUATION
21
22 DX IS THE DISTANCE BETWEEN NEIGHBORING COLUMNS
23
24 DZ IS THE DISTANCE BETWEEN NEIGHBORING ROWS
25
26 IT IS THE MAXIMUM NUMBER OF ITERATION WHICH WILL BE PERFORMED
27
28 WHEN A SUB RETURNS CONTROL TO THE MAIN PROGRAM IT WILL
29 EQUAL THE NUMBER OF ITERATIONS PERFORMED
30
31 ***** THE USE ***** SOLVE POISSON EQUATION *****
32
33 CONSIDER A RECTANGULAR MESH X(I,J) WHICH HAS N ROWS AND M COLUMNS.
34 THE VALUES OF THE BOUNDARY ARE GIVEN. THAT IS THE VALUES OF ROW
35 I= 0 AND N, COLUMN J= 1 AND COLUMN M ARE KNOWN.
36
37 PART TO OBTAIN INTERIOR VALUES SUCH THAT
38  $X(I,J) = \frac{1}{4} [X(I-1,J) + X(I+1,J) + X(I,J-1) + X(I,J+1)] - G \cdot \text{ETAIL}(I,J)$ 
39 WHERE  $A = 2 \cdot Y \cdot DZ \cdot DX = 0.5 \cdot (A + B) \cdot V = \frac{1}{2} \cdot (A + B) \cdot G = W8$ 
40
41 ***** THE METHOD ***** MODIFIED SUCCESSIVE OVER-RELAXATION *****
42 YOUNG'S METHOD (1)
43
44 FIRST DIVIDE THE MESH POINTS INTO TWO SETS. THE FIRST SET CONTAINS
45 ODD POINTS AND THE SECOND SET CONTAINS THE EVEN POINTS. X(I,J) IS
46 ODD IF I+J IS ODD. IF I+J IS EVEN, THEN X(I,J) IS EVEN.
47
48 EACH ITERATION IS DIVIDED INTO TWO PARTS. THE ODD PART AND THE
49 EVEN PART. THE ODD PART IS DONE FIRST. THE ODD FORMULA IS
50  $X(I,J) = \frac{1}{4} [X(I-1,J) + X(I+1,J) + X(I,J-1) + X(I,J+1)] - G \cdot \text{ETAIL}(I,J)$ 
51 (L)
52
53 (L+1)
54
55 (W-1) * X(I,J) WHERE X(I,J) STANDS FOR X(L,J) AFTER L ITERATIONS.
56 THE EVEN FORMULA IS
57 (L+1)
58
59 (L+1)
60
61 X(I,J) =  $\frac{1}{4} [X(I-1,J) + X(I+1,J) + X(I,J-1) + X(I,J+1)] - G \cdot \text{ETAIL}(I,J)$ 
62 (L)
63
64 (W-1) * X(I,J) IN BOTH FORMULAS W IS THE RELAXATION PARAMETER.
65
66 ***** THE ALGORITHM *****
67
68 THIS ALGORITHM HAS TWO PHASES. PHASE1 IMPROVES 64 POINTS AT A TIME
69
70

```

```

51 8 UNTIL ALL THE POINTS IN CLASS ONE ARE IMPROVED. PHASE2 IMPROVES
52 8 ALL THE POINT IN CLASS TWO, TWO ROWS AT A TIME.
53 8 CONSIDER THE MESH AT * J WHERE FOR 1<1<N AND 1<J<M X(1,J) IS TO BE
54 8 IMPROVED. DIVIDE THESE POINTS INTO TWO CLASSES.
55 8 CLASS ONE CONSISTS OF X(I,J) SUCH THAT 1<1<N AND 1<J<M-1 MOD 64
56 8 CLASS TWO CONSISTS OF X(I,J) SUCH THAT 1<1<N AND (M-1-(M-2)MOD 64)<
57 8 J<M. TO IMPLEMENT THIS I USE STRAIGHT ROW WRAPAROUND STORAGE.
58 8
59 8 ***** CODE NAME: MSOP *****
60 8
61 8 ***** FUNCTION: *****
62 8
63 8 SOLVES POISSONS EQUATION
64 8
65 8 ***** LANGUAGE: GLYPHIC *****
66 8
67 8 ***** REFERENCES *****
68 8
69 8
70 8 ([1] SECTION 2.3 OF THIS PAPER
71 8 [2] VARGA, R. S., "MATRIX ITERATIVE ANALYSIS",
72 8 PRENTICE-HALL,ENGLEWOOD CLIFFS, NEW JERSEY(1962).
73 8 [3] LAWRIE, D. H., "GLYPHIC PROGRAMMING MANUAL",
74 8 ILLIAC IV DOCUMENT NO. 232, DEPARTMENT OF COMPUTER
75 8 SCIENCE, UNIVERSITY OF ILLINOIS AT URBANA-CHAMPAIGN,
76 8 URBANA. ILLINOIS(AUGUST, 1970).
77 8
78 8 *****
79 8
80 8 BEGIN
81 8 BOOLEAN MODE,OLDMODE:
82 8 CINT G,CPI:
83 8 CINT I1,CNM2,CNM3,CM1,CMIK,C1,CCN,CODD,CJ,CK:
84 8 CINT C:
85 8 CREAL EPAS,OMEGAL,ALPH ,BETA,G1:
86 8 CREAL W,W1,AL,BE:
87 8 CINT Y:
88 8 PINT NP:
89 8 PINT JN1:
90 8 PINT PJ,J1,JLN,PN:
91 8 PINT PN2,Jp:
92 8 LABEL DONE:
93 8 W:=WB:
94 8 W1:=1-WH:
95 8 CM:=CM-1:
96 8 CM1:=C1,(16:421:
97 8 BE:=DX*DX:
98 8 AL:=DZ*DZ:
99 8 EPAS:=0.5/(3E+AL):
100 8 G1:=W*AL*RE/(2.0*(AL+BE)):
101 8 ALPH:=1.0/BE:
102 8 BETA:=1.0/AL:
103 8 CPI:=CM1+1:
104 8 Y IS THE NUMBER OF ROWS OF PE MEMORY IT TAKES TO STORE ONE ROW OF
105 8 A CN BY CM MESH.
106 8 Y:=CPI:
107 8 G:=2*CPI:
108 8 IF CODD EQUALS ONE WE HAVE AN ODD NUMBER OF ROWS TO IMPROVE.
109 8 CODD:=CN,(63:1):
110 8 CMIK IS USED TO HANDLE CASE WHERE PHASE 2 MUST IMPROVE
111 8 AN ELEMENT IN PE 63.

```

```

111 CMK:=CM*(58:61)
112 IF CMK=0 THEN BEGIN
113   CM1:=CM1-1;CMK:=64;
114   END ELSE IF CMK=1 THEN CMK=0;

00011100
115 CM3:=CN-3;
116 CM3:=CM3*CP1;
117 * PJ*NP*PN*CN AND CNM2 ARE USED TO HELP EACH PE ACCESS THE RIGHT
118 * MEMORY LOCATION, THAT IS OBTAIN THE FOUR NEIGHBORS IN THE FIVE
119 * POINT DIFFERENCE EQUATION.
120 PJ:=0;PN:=0;
121 MODE:=BOOLEAN(8000000000000000(16));
122 PN:=1;
123 MODE:=BOOLEAN(4000000000000000(16));
124 NP:=1;
125 CNM2:=CN-2;
126 CM2:=CM2*CP1;
127 MODE:=BOOLEAN(5555555555555555(16));
128 PJ:=CP1;
129 OLDMODE:=BOOLEAN(7FFFFFFFFFFFFFFF(16));
130 MODE:=TRUE;
131 CNM1:=CM1-1;
132 * THIS LOOP IMPROVES ALL MESH POINTS. THUS EACH TIME IT IS EXECUTED
133 * WE INCREASE THE ITERATION COUNTER BY ONE.
134 LOOP I1:=1;IT DO BEGIN
135 * THE ITERATION IS DIVIDED INTO TWO PARTS
136 * IF THERE ARE M POINTS TO BE IMPROVED IN EACH ROW WHERE M=64N+1.
137 * N AND I ARE INTEGERS AND I<64 THEN PART ONE IMPROVES THE FIRST 64N
138 * POINTS AND PART TWO IMPROVES THE REMAINING I POINTS ON EACH ROW.
139 C:=0;
140 LOOP CK:=1;1,2 DO BEGIN
141 * PART ONE IMPROVES 64 POINTS AT A TIME IN TWO ADJACENT ROWS.
142 LOOP CJ:=0;1;CCH DO BEGIN
143   J1:=P1+CP1;
144   JN1:=I1J1;
145   J1N:=JN1+NP*PN;
146   J1:=J1+PN;
147 * THIS LOOP IMPROVES LINES TWO AT A TIME THUS IF THERE IS AN ODD
148 * NUMBER OF LINES WE MUST DEAL WITH THE LAST LINE IN A SPECIAL WAY.
149   LOOP C1:=CP1*G*CM3 DO BEGIN
150     PN2:=X(J1);
151     X(J1)=G1*(X(J1+Y)*X(J1-Y))*ALPHA+BETA*(RTL(1,*X(J1N))
152     +RTL(1,*X(JN1)))- ETAT(J1))+W1*X(J1);
153     J1:=J1+G;
154     JN1:=JN1+G;
155     J1N:=J1N+G;
156   IF ABS(PN2-X(J1)) GEQ BOUND THEN C:=1;
157   END;

00014800
158 * IF WE HAVE AN ODD NUMBER OF LINES THE LAST LINE IS IMPROVED HERE.
159 IF CCHD=1 THEN BEGIN
160   JN1:=CNM2+CJ;
161   J1:=JN1+PN;
162   J1N:=J1+NP;
163   PN2:=X(J1);
164   X(J1)=G1*(X(J1+Y)+X(J1-Y))*ALPHA+BETA*(RTL(1,*X(J1N))
165   +RTL(1,*X(JN1)))- ETAT(J1))+W1*X(J1);
166   IF ABS(PN2-X(J1)) GEQ BOUND THEN C:=1;
167   END;

00015800

```

```

00011000 02:00:000155
00011100 02:00:000161
00011200 03:01:000178
00011300 03:01:000190

00011100
00011400 02:00:000225
00011500 02:00:000232
00011600 02:00:000244
00011700 02:00:000244
00011800 02:00:000244
00011900 02:00:000244
00012000 02:00:000248
00012100 02:00:000253
00012200 02:00:000255
00012300 02:00:000260
00012400 02:00:000262
00012500 02:00:000269
00012600 02:00:000281
00012700 02:00:000285
00012800 02:00:000290
00012900 02:00:000295
00013000 02:00:000299
00013100 02:00:000309
00013200 02:00:000306
00013300 02:00:000306
00013400 03:00:000330
00013500 03:00:000330
00013600 03:00:000330
00013700 03:00:000330
00013800 03:00:000330
00013900 03:00:000335
00014000 04:00:000346
00014100 04:00:000346
00014200 05:00:000370
00014300 05:00:000375
00014400 05:00:000405
00014500 05:00:000409
00014600 05:00:000412
00014700 05:00:000412
00014800 05:00:000412
00014900 06:00:000453
00015000 06:00:000460
00015100 06:00:000491
00015200 06:00:000525
00015300 06:00:000530
00015400 06:00:000535
00015500 06:00:000540
00015600 06:00:000575

00014800
00015700 05:00:000583
00015800 05:00:000583
00015900 06:01:000600
00016000 06:01:000606
00016100 06:01:000609
00016200 06:01:000612
00016300 06:01:000619
00016400 06:01:000650
00016500 06:01:000684
00016600 06:01:000719

00015800

```

```

168 PJ:=PJ+1;
169 END;

170 JI:=PJ;
171 JN1:=(11J1;
172 * PART TWO IMPROVES THE REMAINING POINTS.
173 IF PENCMK AND OLDMODE THEN BEGIN
174   DMODE:=MODE;
175 * THIS LOOP IMPROVES TWO LINES AT A TIME.
176 LOOP CI:=CPI+6,CNM3 DO BEGIN
177   MODE:=TRUE;
178   JI:=JI+CI;
179   JN1:=JN1+CI;
180   JN1:=JN1+PN;
181   MODE:=DMODE;
182   P2:=X(J1);
183   X(J1):=G1*(X(J1)+Y1+X(J1-Y1))*ALPH+BETA*(RTL(1,X(JN1))
184   +RT(1,X(JN1)))-ETA(J1))+W1*X(J1);
185   IF ABS(P2-X(J1)) GEO BOUND THEN C:=1;
186   END;

187   MODE:=TRUE;
188   JN1:=CNM2+CM1;
189   JI:=JN1+PN;
190   MODE:=DMODE;
191 * THIS LOOP IMPROVES THE LAST LINE IF THERE IS AN ODD NO. OF LINES.
192   IF CMOD=1
193     P2:=X(J1);
194     X(J1):=I*(X(J1)+Y1+X(J1-Y1))*ALPH+BETA*(RTL(1,X(J1))
195     +RT(1,X(JN1)))-ETA(J1))+W1*X(J1);
196     IF ABS(P2-X(J1)) GEO BOUND THEN C:=1;
197     END;

198   END;

199   PJ:=(11P,J-CM1;
200   END;

201   IF C=0 THEN GO TO DONE;
202   END;

203 DONE: IT:=11;
204   END.

```

NUMBER OF ERRORS DETECTED=0000. NUMBER OF WARNINGS NOTED=0000.  
 SOURCE PROGRAM SIZE= 00204 CARD IMAGES, 001637 SYNTACTIC ITEMS.  
 SOURCE FILE: CARD=ERICKSEN/PRINT/MSOR.  
 MACRO LIBRARY: ILLIAC/GLYPHIR3/MACROS.  
 ESTIMATED CORE MEMORY REQUIREMENTS= 000717 WORDS CODE, 000868 WORDS DATA.  
 COMPIATION TIME= 0018 SECONDS ELAPSED, 0009 SECONDS PROCESSING.



## APPENDIX D

## SUCCESSIVE LINE OVER-RELAXATION METHOD IN GLYPNIR, SLOR

## ILLIAC IV GLYPNIR COMPILER

( 86500 VERSION 3.3.00 )

```

1 SUBROUTINE MSLOR(PCPOINT OUT U,CREAL W,CINT CN,CINT CM,CREAL
2 BOUND,PCPOINT I,CREAL DX,CREAL DZ,CINT OUT IT);
3
4 ***** DEFINITION OF PARAMETERS *****
5
6 U IS THE MESH TO BE EVALUATED IN POISSON'S EQUATION
7 T IS THE CONSTANT MESH
8 W IS THE RELAXATION PARAMETER WHICH IS SUPPLIED BY THE
9 MAIN PROGRAM
10 CM IS THE NUMBER OF COLUMNS IN U
11 CN IS THE NUMBER OF ROWS IN U
12 DX IS THE DISTANCE BETWEEN ROWS IN U
13 DY IS THE DISTANCE BETWEEN COLUMNS IN U
14 IT ENTERS AS THE MAXIMUM NUMBER OF ITERATIONS TO BE
15 PERFORMED AND IS RETURNED TO THE MAIN PROGRAM AS THE
16 NUMBER OF ITERATIONS PERFORMED
17
18 ***** THE USE ***** SOLVES POISSON'S EQUATIONS
19
20 CONSIDER A RECTANGULAR MESH, U(I,J), WHICH HAS CN ROWS AND
21 CM COLUMNS. THE VALUES OF THE BOUNDARY ARE GIVEN. THAT IS
22 THE VALUES OF ROW 1, ROW CN, COLUMN 1, AND COLUMN CM ARE
23 KNOWN. THE CONSTANT MESH T IS ALSO KNOWN. WE WANT TO OBTAIN
24 THE INTERIOR POINTS OF U SUCH THAT:
25  $U(I,J) = W(U(I-1,J) + U(I+1,J)) + W(U(I,J-1) + U(I,J+1)) - G * T(I,J)$ 
26 WHERE  $A = DY * DY$ ,  $B = DX * DX$ ,  $H = A / (2 * (A+B))$ ,  $V = B / (2 * (A+B))$ , AND  $G = H * B$ 
27 THIS PROGRAM IS WRITTEN TO WORK ON ANY RECTANGULAR MESH
28 THAT FITS ON CORE.
29
30 ***** METHOD ***** SUCCESSIVE LINE OVERRELAXATION *****
31
32 THIS PROGRAM USES THE CUTHILL-VARGA METHOD (3) PAGE 238 TO
33 PERFORM COLUMN RELAXATION ON THE MESH U WHICH IS STORED IN
34 ODD-EVEN STORAGE AS FOUND IN (1) PAGE 22. T IS ALSO IN
35 ODD-EVEN STORAGE.
36 THE COEFFICIENTS CALCULATED IN PRECONDITIONING, (2) PAGE 32, ARE
37 DISB,DI,D, AND E.
38
39 ***** CODE NAME: MSLOR *****
40
41 ***** FUNCTION: *****
42
43 ***** SOLVES POISSON'S EQUATION *****
44
45 ***** LANGUAGE: GLYPNIR *****
46
47 ***** REFERENCES *****
48
49 (1) SECTION 2.3.4 OF THIS PAPER
50 (2) SECTION 2.4.4 OF THIS PAPER

```

```

00000100 00:00:000000
00000200 01:00:000012
00000300 01:00:000012
00000400 01:00:000012
00000500 01:00:000012
00000600 01:00:000012
00000700 01:00:000012
00000800 01:00:000012
00000900 01:00:000012
00001000 01:00:000012
00001100 01:00:000012
00001200 01:00:000012
00001300 01:00:000012
00001400 01:00:000012
00001500 01:00:000012
00001600 01:00:000012
00001700 01:00:000012
00001800 01:00:000012
00001900 01:00:000012
00002000 01:00:000012
00002100 01:00:000012
00002200 01:00:000012
00002300 01:00:000012
00002400 01:00:000012
00002500 01:00:000012
00002600 01:00:000012
00002700 01:00:000012
00002800 01:00:000012
00002900 01:00:000012
00003000 01:00:000012
00003100 01:00:000012
00003200 01:00:000012
00003300 01:00:000012
00003400 01:00:000012
00003500 01:00:000012
00003600 01:00:000012
00003700 01:00:000012
00003800 01:00:000012
00003900 01:00:000012
00004000 01:00:000012
00004100 01:00:000012
00004200 01:00:000012
00004300 01:00:000012
00004400 01:00:000012
00004500 01:00:000012
00004600 01:00:000012
00004700 01:00:000012
00004800 01:00:000012
00004900 01:00:000012
00005000 01:00:000012

```

```

51 * [3] CUTHILL AND VARGA, "A METHOD OF NORMALIZING BLOCK
52 * ITERATION", J. ASSOC. COMPUT. MACH., VOL. 6, PP. 236-
53 * 244, 1959.
54 * [4] LAWRIE, D. H., "GLYPNIR PROGRAMMING MANUAL",
55 * ILLIAC IV DOCUMENT NO. 232, DEPARTMENT OF COMPUTER
56 * SCIENCE, UNIVERSITY OF ILLINOIS AT URBANA-CHAMPAIGN,
57 * URBANA, ILLINOIS, 1970.
58 *
59 * *****
60 *
61 BEGIN
62 PCPOINT DISB;
63 CINT MC;
64 PCPOINT STORE,E,D,DI;
65 CINT CM2,CM3,CM2K,OD,C,B,EV,CM4,LPR,CL,CJ,IL,CP,LP,ODD,CN2;
66 PINT PN,LPN;
67 PREAL ST;
68 CINT L,K,KPKM1,JC;
69 CREAL ALPHA,RABONE,9FIA,GAMMA,ALPHS;
70 BOUNLEAN OLMODE,DMODE;
71 LABEL DONE;
72 CM2:=CM-2; %THE NUMBER OF COLUMNS TO BE IMPROVED
73 STORE:=GETPEB(CM2); % TEMPORARY STORAGE NEEDED FOR OVERRELAXATION
74 LPR:=CM, (16:42) % LINES OF PEM FOR STORING D,DI OR E MINUS ONE.
75 IF CM, (58:6) GEJ 3 THEN LPR:=LPR+1;
76 D:=GETPEB(LPR); %DIAGONAL CONSTANTS USED BY CUTHILL + VARGA.
77 DISH:=GETPEB(LPR);
78 DI:=GETPEB(LPR); % INVERSE OF DI
79 E:=GETPEB(LPR); % OTHER CONSTANTS IN CUTHILL-VARGA METHOD.
80 LPR:=LPR-1;
81 ALPH:= -1.0/(DZ*IDZ); % ROW ELEMENTS OFF THE DIAGONAL
82 HFTA:= -1.0/(DX*IDX); % COLUMN ELEMENTS OFF THE DIAGONAL
83 GAMMA:= -2.0*(ALPH+BETA); %THE DIAGONAL ELEMENTS
84 ALPHS:=ALPH*ALPH;
85 DMODE:=BOOLEAN(8000000000000000(16));
86 OLMODE:=BOOLEAN(7FFFFFFFFFFFFFFF(16));
87 CN2:=CN-2; % NUMBER OF ROWS TO BE IMPROVED
88 CN2K:=CM2, (57:6); % NUMBER OF PES USED IN PHASE TWO
89 OD:=CN2, (16:41); % (# OF SETS OF 64 ODD ROWS IMPROVED BY PHASE 1
90 IF CN, (57:7)=1 THEN K:=3 ELSE K:=2;
91 K:=K+OD*OD;
92 EV:=(CN-1), (16:41); % # OF SETS OF 64 EVEN ROWS PHASE 1 IMPROVED
93 KPKM1:=K*K-1;
94 CM3:=CM-3;
95 CM4:=CM-4;
96 PN:=0; % VARIABLE USED AS AN INDEX
97 MODE:=BOOLEAN(1);
98 PN:=1;
99 MODE:=TRUE;
100 IF PEN=0 THEN D(0):=GAMMA;
101 * THE FOLLOWING SECTION CALCULATES D,DI AND E
102 LOOP CL:=1, CM3 DO IF CL, (58:6)=PEN THEN BEGIN
103 C:=CL, (16:42); D(C):=GAMMA-ALPHS/ RTR(1, D(C-PN)); END;
104 B:=CM2, (58:6);
105 IF B NEQ 0 AND PEN GEQ 8 THEN D(LPR):=1.0;
106 LOOP CL:=0, 1, LPR DO BEGIN
107 D(C1):=SORT(D(C1));
108 D(C1):=1.0/D(C1);
109 DISH(C1):=D(C1)*O1(C1)*BETA;

```

00010200

```

00010400
00010500
00010600
00010700
00010800
00010900
03:01:000356
03:00:000401
03:00:000407
03:00:000453
03:00:000477
03:00:000488
03:00:000501

```

```

00005100
00005200
00005300
00005400
00005500
00005600
00005700
00005800
00005900
00006000
00006100
00006200
00006300
00006400
00006500
00006600
00006700
00006800
00006900
00007000
00007100
00007200
00007300
00007400
00007500
00007600
00007700
00007800
00007900
00008000
00008100
00008200
00008300
00008400
00008500
00008600
00008700
00008800
00008900
00009000
00009100
00009200
00009300
00009400
00009500
00009600
00009700
00009800
00009900
00010000
00010100
00010200
00010300

```



```

110 E(CI):=ALPH*DI(CI):END;
00010600 00011000 03:00:000517

111 MODE:=OLD MODE;
112 E(0):=F(0)+RTR(1,DI(0));
113 MODE:=TRUE;
114 LOOP CI:=1,LP DO E(CI):=E(CI)*RTR(1,DI(CI-PN));
115 ODD:=CN.[63:1];% IF ODD=1 THEN ROWS TO BE IMPROVED IS ODD.
116 % PREPARE FIRST AND LAST COLUMN OF MESH FOR ITERATION
117 LPN:=K-1;
118 IF PEN=0 THEN LPN:=0;
119 LOOP CI:=0,1,CM3 DO
120 BEGIN LPN:=LPN+K;
121 MC:=CI.[16:42];
122 RABONE:=GRABONE(D(MC),CI);
123 IF CN2K=0 THEN U(LPN+K-1):=RABONE*U(LPN+K-1);
124 IF PEN=0 OR PEN=CN2K THEN
125 U(LPN):=RABONE*U(LPN):END;

126 LP:=0;CJ:=CM-1;
127 % ADD THE FIRST AND THE LAST ROW OF THE MESH TO T
128 C:=CM3*K;
129 LOOP CI:=K,1,KPM DO BEGIN
130 T(CI):=T(CI)+ALPH*U(CI-K);
131 CP:=CI+C;
132 T(CP):=T(CP)+ALPH*U(CP-K);
133 END;

134 % MULTIPLY T BY D INVERSE, DI
135 CP:=K;
136 LOOP CI:=0,1,CM3 DO LOOP CJ:=0,1,K-1 DO BEGIN
137 II:=CI.[16:42];
138 T(CP):=-T(CP)+GRABONE(D(II),CI);
139 CP:=CP+1;END;

140 LOOP II:=1,1,IT DO BEGIN % EACH TIME THIS LOOP CYCLES IS ONE ITER.
141 B:=0;
142 % PHASE ONE STARTS HERE
143 LOOP CI:=1,1,EV DO BEGIN
144 % LPN ACCESSES THE ODD COLUMNS, LP ACCESSES THE EVEN COLUMNS
145 % THIS PART OF PHASE 1 IMPROVES THE EVEN COLUMNS
146 LP:=CI+K;
147 LPN:=LP-1;
148 MODE:=DMODE;LPN:=LPN+2; MODE:=TRUE;
149 STORE(IJ):=T(LP)-GRABONE(DIS(0),0)*(RTL(1,JI(LPN))+U(LP-1));
150 LOOP CJ:=1,1,CM3 DO BEGIN C:=CJ.[16:42];
151 LP:=LP+K;LPN:=LPN+K;
152 STORE(CJ):=T(LP)-GRABONE(DIS(C),CJ)*(RTL(1,JI(LPN))+U(LP-1))-
153 GRABONE(E(C),CJ)*STORE(CJ-1);
154 END;

155 LOOP MC:=1,1,CM3 DO BEGIN
156 CJ:=CM3-MC;
157 JC:=CJ+1;
158 C:=JC.[16:42];
159 STORE(CJ):=STORE(CJ)-GRABONE(E(C),JC)*STORE(JC);
160 END;

161 LP:=CI+K;
162 LOOP CJ:=0,1,CM3 DO BEGIN % PERFORM OVERRELAXATION
163 ST:=U(LP1);

```

```

164 U(LP):=W*(STORE(CJ)-U(LP))+U(LP);
165 IF ABS(U(LP)-ST)>BOUND THEN B:=1; % CHECK ERROR
166 LP:=LP*K;END;
167
168
169 PHASE TWO STARTS HERE
170 IF CN2K+ODD>0 AND CN2K+ODD<64 THEN BEGIN
171 LP:=KPKM1;
172 LPN:=LP-1;
173 STORE(0):=T(LP1-GRABONE(D1SB(0),0))*(RTL(1,,U(LP1))+U(LP-1));
174 LOOP CJ:=1,1,CM3 DO BEGIN C:=CJ.{16:42};
175 LP:=LP*K;LPN:=LPN*K;
176 STORE(CJ):=T(LP1-GRABONE(D1SB(C),CJ))*(RTL(1,,U(LP1))+U(LP-1))-
177 GRABONE(E(C),CJ)*STORE(CJ-1);
178 END;
179
180 LOOP MC:=1,1,CM3 DO BEGIN
181 CJ:=CM3-MC;
182 JC:=CJ+1;
183 C:=JC.{16:42};
184 STORE(CJ):=STORE(CJ)-GRABONE(E(C),JC)*STORE(JC1);
185 END;
186
187 LP:=KPKM1;
188 LOOP CJ:=0,1,CM3 DO BEGIN % PERFORM OVERRELAXATION
189 STORE(CJ):=W*(STORE(CJ)-U(LP))+U(LP);
190 IF (ODD+CN2K) > PEN THEN BEGIN
191 IF ABS(U(LP1)-STORE(CJ))>BOUND THEN B:=1; % CHECK ERROR
192 U(LP):=STORE(CJ);END;
193 JC:=CJ+1;
194 LP:=LP*K;END;
195
196
197
198
199 PHASE ONE STARTS HERE
200 LNP WILL BE USED TO ACCESS THE ODD COLUMNS, LP THE EVEN
201 LOOP CI:=1,1,OD DO BEGIN
202 LP:=CI*K;
203 LPN:=LP-1;
204 MODE:=DMODE;LPN:=LPN+2; MODE:=TRUE;
205 STORE(0):=T(LP1-GRABONE(D1SB(0),0))*(RTL(1,,U(LP))+U(LP1));
206 LOOP CJ:=1,1,CM3 DO BEGIN C:=CJ.{16:42};
207 LP:=LP*K;LPN:=LPN*K;
208 STORE(CJ):=T(LP1-GRABONE(D1SB(C),CJ))*(RTL(1,,U(LP))+U(LP1))-
209 GRABONE(E(C),CJ)*STORE(CJ-1);
210 END;
211
212 LOOP MC:=1,1,CM3 DO BEGIN
213 CJ:=CM3-MC;
214 JC:=CJ+1;
215 C:=JC.{16:42};
216 STORE(CJ):=STORE(CJ)-GRABONE(E(C),JC)*STORE(JC);
217 END;
218
219 LPN:=K+C1-1;
220 MODE:=DMODE;LPN:=LPN+2; MODE:=TRUE;
221 LOOP CJ:=0,1,CM3 DO BEGIN % PERFORM OVERRELAXATION
222 ST:=U(LP1);

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00016400 05:00:001347
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00020300 05:00:002015
00020400 05:00:002040
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00020600 05:00:002072
00020700 05:00:002080
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00021400 05:00:002170

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215 U[LPN]:=W*(STORE[CJ]-U[LPN])+U[LPN];
216 IF ABS(U[LPN]-ST)>BOUND THEN B:=1; % CHECK ERROR
217 LPN:=LPN+K;END;
218 END;
219 *
220 PHASE TWO STARTS HERE
221 IF CN2K>0 THEN BEGIN
222 LP:=KPKM1;
223 LPN:=LP-1;
224 STORE[0 ]:=T[LPN]-GRABONE(D[ISB(0),0])*(RTR(1,,U[LP])+U[LP]);
225 LP:=LP+K;LPN:=LPN+K;
226 STORE[CJ]:=T[LPN]-GRABONE(D[ISB(CJ),CJ])*(RTR(1,,U[LP])+U[LP])-
227 GRABONE(E[CJ],CJ)*STORE[CJ-1];
228 END;
229
230 LOOP MC:=1,1,CM3 DO BEGIN
231 CJ:=CM3-MC;
232 JC:=CJ+1;
233 C:=JC.[16:42];
234 STORE[CJ]:=STORE[CJ]-GRABONE(E[CJ],JC)*STORE[JC];
235 END;
236
237 LPN:=K+K-2;
238 LOOP CJ:=0,1,CM3 DO BEGIN % PERFORM OVERRELAXATION
239 STORE[CJ]:=W*(STORE[CJ]-U[LPN])+U[LPN];
240 IF OLDMODE AND CN2K GEQ PEN THEN BEGIN
241 IF ABS(U[LPN]- STORE[CJ])>BOUND THEN B:=1; % CHECK ERROR
242 U[LPN]:=STORE[CJ];END;
243 LPN:=LPN+K;END;
244
245 END;
246
247 IF B=0 THEN GO TO DONE;END;
248
249 DONE: I1:=I1;
250 CP:=K;
251 LOOP CI:=0,1,CM3 DO LOOP CJ:=0,1,K-1 DO BEGIN
252 I1:=CI.[16:42];
253 U[CP]:= U[CP]*GRABONE(D[I1],CI);
254 T[CP]:=-T[CP]*GRABONE(D[I1],CI);
255 CP:=CP+1;END;
256
257 C:=K*CM3;
258 LOOP CI:=K,1,KPKM1 DO BEGIN
259 T[CI]:=T[CI]-ALPH*U[CI-K];
260 CP:=CI+C;
261 T[CP]:=T[CP]-ALPH*U[CP+K];END;
262
263 END.*% OF MSLOR
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NUMBER OF ERRORS DETECTED=0000. NUMBER OF WARNINGS NOTED=0000.  
SOURCE PROGRAM SIZE= 00256 CARD IMAGES, 003186 SYNTACTIC ITEMS.  
SOURCE FILE: CARD=ERICKSEN/PRINT/SLOR.

MACPO LIBRARY: ILLIAC/GLYPNIR3/MACROS.  
ESTIMATED CORE MEMORY REQUIREMENT= 001707 WORDS CODE, 000806 WORDS DATA.  
COMPILATION TIME= 0034 SECONDS ELAPSED, 0017 SECONDS PROCESSING.

APPENDIX E

PEACEMAN-RACHFORD ADI IN GLYPNIR  
ILLIAC IV GLYPNIR COMPILER  
( B6500 VERSION 3.3.00 )

```

1 SUBROUTINE ADI(PCPOINT OUT U,PCPOINT I,CINT CM,CINT CN,CREAL HX,
2 CREAL HY,CINT FLAG,CINT OUT MAXIT,CREAL ERROR);
3
4 %%%%%%%%%% DEFINITION OF PARAMETERS %%%%%%%%%%
5
6 U IS THE MESH TO BE IMPROVED
7 T IS THE CONSTANT MESH IN POISSON'S EQUATION
8 CM IS THE NUMBER OF COLUMNS IN U
9 CN IS THE NUMBER OF ROWS IN U
10 HX IS THE DISTANCE BETWEEN ROWS IN U
11 HY IS THE DISTANCE BETWEEN COLUMNS IN U
12 FLAG TELLS THE SUBROUTINE IF THE MAIN PROGRAM SUPPLIED
13 THE NUMBER OF ITERATIONS TO BE PERFORMED OR THE
14 ERROR BOUND
15 MAXIT IS THE NUMBER OF ITERATIONS THAT WILL BE PERFORMED
16 IF FLAG=0 THEN MAXIT MUST BE SUPPLIED BY THE MAIN PROGRAM
17 ERROR IS THE ERROR BOUND OBTAINED AFTER MAXIT ITERATIONS
18 IF FLAG=1 THEN THE MAIN PROGRAM SUPPLIES ERROR AND MAXIT IS
19 CALCULATED IN THIS SUBROUTINE
20
21 %%%%%%%%%% THE USE %%%%%%%%%% SOLVE POISSON'S EQUATION %%%%%%%%%%
22
23 CONSIDER A RECTANGULAR MESH, U(I,J), WHICH HAS CN ROWS AND
24 CM COLUMNS. THE VALUES OF THE BOUNDARY ARE GIVEN. THAT
25 IS THE VALUES OF ROW 1, ROW N, COLUMN 1 AND COLUMN M ARE
26 KNOWN. THE CONSTANT MESH T IS ALSO KNOWN. WE WANT TO
27 OBTAIN THE INTERIOR VALUES OF U SUCH THAT:
28  $U(I,J) = H * (U(I-1,J) + U(I+1,J)) + V * (U(I,J-1) + U(I,J+1)) - G * T(I,J)$ 
29 WHERE  $A = H * H * Y$ ,  $B = H * H * X$ ,  $H = A / (2 * (A * B))$ ,  $V = B / (2 * (A * B))$ , AND  $G = H * B$ 
30 THIS PROGRAM IS LIMITED TO MESHES WHERE CN AND CM ARE
31 LESS THAN OR EQUAL TO 64.
32
33 %%%%%%%%%% METHOD %%%%%%%%%% ALTERNATING DIRECTION IMPLICIT METHOD %
34
35 WE WILL USE THE PEACEMAN-RACHFORD SCHEME AS FOUND IN [2] PAGE
36 174. WE CALCULATE THE OPTIMAL RELAXATION PARAMETERS FOR ALL THE
37 ITERATIONS USING METHOD IN [2] PAGE 192. THEN WE DO THE
38 PRECONDITIONING FOR THOMAS' METHOD OF INVERTING TRIANGULAR
39 MATRICES AS FOUND IN [1] PAGE 37.
40 FINALLY WE DO THE ROW RELAXATION FOLLOWED BY THE COLUMN
41 RELAXATION.
42
43 %%%%%%%%%% TYPE OF STORAGE USED FOR THE MESHES %%%%%%%%%%
44
45 U AND T ARE TO BE SUPPLIED TO THE SUBROUTINE IN STRAIGHT
46 STORAGE. BEFORE ANY ITERATIONS ARE PERFORMED U AND T WILL
47 WILL BE PLACED IN SKEWED STORAGE. AFTER THE LAST ITERATION
48 U AND T WILL BE RETURNED TO STRAIGHT STORAGE.
49
50 %%%%%%%%%% LANGUAGE GLYPNIR %%%%%%%%%%

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00000100 00:00:000000
0000200 01:00:000012
0000300 01:00:000012
0000400 01:00:000012
0000500 01:00:000012
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0001200 01:00:000012
0001300 01:00:000012
0001400 01:00:000012
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0001900 01:00:000012
0002000 01:00:000012
0002100 01:00:000012
0002200 01:00:000012
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0002400 01:00:000012
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0004600 01:00:000012
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51	%	REFERENCES	%	0005100
52	%	SECTION 2.5 OF THIS PAPER	%	00005200
53	%	WACHPRESS, E. L., "ITERATIVE SOLUTION OF ELLIPTIC	%	00005300
54	%	SYSTEMS AND APPLICATIONS TO NEUTRON DIFFUSION EQUATIONS	%	00005400
55	%	OF REACTOR PHYSICS", RENTICE-HALL, ENGLEWOOD CLIFFS,	%	00005500
56	%	NEW JERSEY, 1966.	%	00005600
57	%	LAWRIE, D. H., "GLYPNIR PROGRAMMING MANUAL".	%	00005700
58	%	ILLIAC IV DOCUMENT NO. 232. DEPARTMENT OF COMPUTER	%	00005800
59	%	SCIENCE, UNIVERSITY OF ILLINOIS AT URBANA-CHAMPAIGN,	%	00005900
60	%	URBANA, ILLINOIS, 1970.	%	00006000
61	%		%	00006100
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64	%		%	00006400
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165 CA:=2.0*BETA*GRABONE(W,CC);
166 LOOP CI:=1,CN2 DO BEGIN
167   MODE:=TRUE;RABONE:=GRABONE(SICI-1,CC);
168   HMODE:=REVR(1,HMODE);
169   MODE:=HMODE;
170   U(CI):=(AI*(-T(CI)-BETA*(RTR(1,U(CI))+RTL(1,U(CI)))+CA*U(CI))
171   -RTR(1,U(CI-1)))*RABONE;END;

172 LOOP IC:=0,1,CN3 DO BEGIN
173   CI:=CN2-IC;
174   MODE:=TRUE;RABONE:=GRABONE(SICI-1,CC);
175   HMODE:=HMODE;
176   U(CI):=U(CI)-RABONE*RTL(1,U(CI+1));
177   HMODE:=REVL(1,HMODE);END;

178   MODE:=TRUE;
179   END;

180   CK:=CK+ITER;
181   ITE:=62;
182   ITER:=32;END;

183   LOOP CI:=1,CN-1 DO BEGIN
184     U(CI):=RTL(CI,U(CI));
185     T(CI):=RTL(CI,T(CI));END;

186   END.

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00021800 00021800 02:00:001838  
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NUMBER OF ERRORS DETECTED=0000. NUMBER OF WARNINGS NOTED=0000.  
SOURCE PROGRAM SIZE= 00186 CARD IMAGES, 002145 SYNTACTIC ITEMS.  
SOURCE FILE: CARD=ERICKSEN/PRINT/ADI.  
MACRO LIBRARY: ILLIAC/GLYPNIR3/MACROS.  
ESTIMATED CORE MEMORY REQUIREMENTS= 001084 WORDS CODE, 000684 WORDS DATA.  
COMPILATION TIME= 0037 SECONDS ELAPSED, 0015 SECONDS PROCESSING.

HOCKNEY'S DIRECT METHOD, MODIFIED, MFACR  
ILLIAC IV GLYPHIC COMPILER  
( 86500 VERSION 3.3.00 )

```

1 BEGIN
2   PE INTEGER SUBROUTINE TURN AS RGA(CINT P, PINT I);
3   BEGIN
4     RGA := I; CARL := P;
5     CODE BEGIN
6       LIT(2)      = 1.0, 1;
7       SHABR      0(1);
8       CSHL(1)    24;
9       CLRA;
10      COR(1)     SC2;
11      SHAPL      1;
12      RTAR       2;
13      TREFAM(1)  -3;
14      CSHR(1)    24;
15      RTAL       1(1);
16      END CODE;
17      END; % OF TURN %%%%%%%%%%
18 %%%%%%%%%%
19 SUBROUTINE SETPT1(PCPOINT A, PCPOINT S, PCPOINT ISS, CINT OUT PP
20   , PCPOINT INDEX, CINT NX, CREAL HX, CREAL HY) ;
21   % SETPT1 CALCULATES FOUR PARAMETERS REQUIRED BY
22   % POISSONDIRECT, A+S, ISS, INDEX. THESE PARAMETERS
23   % ARE VECTORS WHICH REQUIRE CN/128 LINES OF PEM
24   % EACH. CN IS THE NUMBER OF ROWS IN THE MESH, NOT
25   % PEM ROWS. NX EQUALS CN-1. HX AND HY ARE
26   % THE HORIZONTAL AND VERTICAL COEFFICIENTS
27   % RESPECTIVELY. PP IS THE LOG BASE 2 OF NX.
28   % SETPT1 MUST BE CALLED EVERY TIME THE MESH SIZE CHANGES
29   % OR HX AND/OR HY CHANGE.
30   BEGIN
31     CINT I, R, KN;
32     PINT L, Q;
33     CREAL P1, H, X;
34     P1 := 3.1415926535898;
35     PP := LN(NX)/LN(2);
36     L := PEN;
37     H := LN(2); SIMWRITE(LINE, "LN(2)", H);
38     X := P1 / NX;
39     H := HX / HY;
40     KN := NX * (16:41);
41     LOOP I := 0, 1, KN DO
42       BEGIN
43         S(I) := SIN(L * X);
44         ISS(I) := S(I) + S(I);
45         IF ISS(I) NEQ 0 THEN BEGIN ISS(I) := -1.0 / ISS(I); END;
46
47         A(I) := -2.0 * (H - H * COS(L * X) + 1);
48         INDEX(I) := TURN(PP-1, L);
49         L := L + 64;

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49      END ;
50      END ;
51      SUBROUTINE
52      POISSONDIRECT(PCPOINT P,PCPOINT C,PCPOINT A,PCPOINT S,CINT CN,
53      CINT CM,CREAL HX,CREAL HY,PCPOINT ISS,
54      CINT LLN,PCPOINT INDEX ) ;
55      ***** FUNCTION *****
56      *****
57      THIS SUBROUTINE SOLVES POISSON' EQUATION USING A DIRECT
58      METHOD SIMILAR TO HOCKNEY'S PROGRAM (1).
59      C(I,J) = HX*(P(I+1,J)+P(I-1,J)-2*P(I,J)) +
60      HY*(P(I,J+1)+P(I,J-1)-2*P(I,J)).
61      THE (CHARGE, C, AND THE BOUNDARY POINTS OF THE POTENTIAL,
62      P, ARE USED TO CALCULATE THE INTERIOR POINTS OF POTENTIAL.
63      *****
64      ***** PARAMETERS *****
65      *****
66      P IS THE POTENTIAL. C IS THE CHARGE.
67      A CONTAINS THE DIAGONAL ELEMENTS OF THE TRIDIAGONAL
68      MATRICES SOLVED IN THE SECTION OF THIS ROUTINE CALLED
69      CRED. S CONTAINS THE VALUES OF THE SIN WHICH ARE NEEDED
70      BY THE FOURIER TRANSFORM PART OF THE ALGORITHM. ISS IS
71      ALSO NEEDED BY FOURIER.
72      INDEX IS USED FOR SCRAMBLING BY FOURIER ANALYSIS AND
73      SYNTHESIS. A, S, ISS, AND INDEX ARE CALCULATED BY SETPT1.
74      P HAS CN ROWS AND CM COLUMNS. HX(HY) IS THE HORIZONTAL(
75      VERTICAL) COEFFICIENT OF POISSON'S EQUATION. LLN EQUALS
76      THE LOG BASE 2 OF CN-1.
77      *****
78      ***** METHOD *****
79      *****
80      A FOURIER ANALYSIS IS DONE TO EACH COLUMN BY SUBROUTINE
81      FOURIER. THIS IS FOLLOWED BY SOLVING THE TRIDIAGONAL
82      MATRICES, ONE PER ROW. FINALLY A FOURIER SYNTHESIS IS
83      PERFORMED ON EACH ROW TO GIVE ANSWER. FOR FURTHER DETAIL
84      SEE SECTIONS 3, 4 AND 5 OF THIS PAPER.
85      *****
86      ***** REFERENCES *****
87      *****
88      (1) R.W. HOCKNEY, "THE POTENTIAL CALCULATION AND
89      SOME APPLICATIONS", METHODS IN COMPUTATIONAL
90      PHYSICS, VOL. 9, P. 136-211, 1970.
91      (2) J. W. COOLEY, ET AL, "THE FAST FOURIER TRANSFORM
92      ALGORITHM: PROGRAMMING CONSIDERATIONS IN THE
93      CALCULATION OF SINE, COSINE AND LAPLACE TRANSFORMS"
94      JOURNAL OF SOUND AND VIBRATION, VOLUME 12 NUMBER 2,
95      JUNE 1970.
96      BEGIN
97      CINT CM2,CN2,K,CNIK,CJ,CL,IL,LL1,L2,L3,L4,KM1,KP,KC,I,
98      L,OFF,SL,I1,J,N2,ILJ,KM,N3,I1,IP;
99      LABEL FOURIER,FINISH ;
100     PINT P1,PJ;
101     CREAL POTFAC,RR,RR1,CO,S1; BOOLEAN FMODE,SMODE,SAVEMODE ;
102     PCPOINT STORE ;
103     PREAL PR,ODD1,ODD2,ODD3 ;
104     PREAL E,O,I;
105     ***** SECTION 1 *****
106     SAVEMODE:=MODE; MODE:=TRUE;

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00004200 00004900 04:00:000280
00004200 00005000 03:00:000288
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00010600 03:00:000306

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107 K := CM .(116:42) ;
108 I := 0 ;
109 IF CM.(58:6) GTR 0 THEN I := 1 ;
110 CN2 := CN - 2 ; POTFAC := 0.125 / ((CN-1)*HY) ;
111 CM2 := CM - 2 ;
112 K := K + I ;
113 KM1 := K - 1 ;
114 KP := K + K ;
115 STORE := GETPEB(KP+K) ;
116 KC := K * CN2 ; PI := 0 ;
117 I := CM.(58:6) ;
118 OFF := 1 - 2 ;
119 IF I = 0 THEN BEGIN OFF := 62 ; END ;

120 MODE := ( PEN GTR OFF ) ;
121 PI := -1 ;
122 MODE := TRUE ;
123 CJ := KC + 1 ;
124 LOOP I := 0+1,KM1 DO
125 BEGIN
126 STORE[I] := C[I+K] ;
127 STORE[I+K] := C[I+KC] ;
128 STORE[I*KP] := P[I] ;
129 END ;

130 LOOP I := 1+1,KM1 DO
131 BEGIN
132 C[PI+CJ] := C[PI+CJ] - P[CJ+PI*K] * HX ;
133 CJ := CJ + 1 ;
134 C[PI+K+I] := C[PI+K+I] - P[PI+I] * HX ;
135 END ;

136 IF PEN GTR 0 AND PEN LSS OFF+1 THEN
137 BEGIN
138 C[K] := C[K] - P[0] * HX ;
139 C[KC] := C[KC] - P[KC*K] * HX ;
140 FMODE := MODE ;
141 END ;

142 PJ := 0 ;
143 PI := 0 ;
144 MODE := ( PEN EQL OFF ) ;
145 PI := KM1 ;
146 MODE := TRUE ;
147 MODE := ( PEN GTR OFF-1 ) ;
148 PJ := -1 ;
149 MODE := TRUE ;
150 IF OFF = 1 THEN BEGIN FMODE := (PEN GTR 0) ; END ;

151 LOOP I := K*K,KC DO
152 BEGIN
153 PR := 0.0 ;
154 MODE := ( PEN EQL OFF ) ;
155 PR := RTL(1,P[I+KM1]) ;
156 MODE := TRUE ;
157 MODE := PEN EQL 1 ;
158 PR := RTL(1,P[I]) + PR ;
159 MODE := TRUE ;
160 MODE := FMODE ;
161 P[1+PI] := (C[I+PI] -PR*HY) * POTFAC ;

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00010700 03:00:000312
00010800 03:00:000318
00010900 03:00:000323
00011000 03:00:000349
00011100 03:00:000380
00011200 03:00:000387
00011300 03:00:000396
00011400 03:00:000403
00011500 03:00:000412
00011600 03:00:000423
00011700 03:00:000437
00011800 03:00:000443
00011900 03:00:000450

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00011900
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00012000 03:00:000473
00012100 03:00:000486
00012200 03:00:000492
00012300 03:00:000496
00012400 03:00:000503
00012500 03:00:000527
00012600 04:00:000527
00012700 04:00:000540
00012800 04:00:000559
00012900 04:00:000573

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00012500
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00013000 03:00:000581
00013100 03:00:000605
00013200 04:00:000605
00013300 04:00:000639
00013400 04:00:000646
00013500 04:00:000680

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00013600 03:00:000688
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00013800 04:01:000717
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00014100 04:01:000762

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00014200 03:00:000764
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00014700 03:00:000788
00014800 03:00:000802
00014900 03:00:000808
00015000 03:00:000812

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00015100 03:00:000841
00015200 03:00:000882
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00015600 04:00:000915
00015700 04:00:000919
00015800 04:00:000929
00015900 04:00:000941
00016000 04:00:000945
00016100 04:00:000949

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162 MODE := TRUE ;
163 LOOP L1:= 1,I,KM1 DO
164 BEGIN
165 P(L1+1+PJ) := C(L1+I+PJ) * POTFAC ;
166 END;
167
168 END ;
169
170 J := (CN-1)*K;
171 L:=J.(16:47);
172 II:= 0;
173 %%%%%%%%%% SECTION 2 %%%%%%%%%%
174 % FOURIER ANALYSIS AND SYNTHESIS
175 FOURIER: II := II + 1;
176 P1 :=0;
177 IF ( PEN EQL OFF+1 ) THEN BEGIN
178 PI := -1 ; END ;
179
180 MODE := FMODE ;
181 KM1 := K-1 ;
182 KP := K + K ;
183 N3 := CN - 1 ;
184 N2 := N3.(16:47);
185 LOOP ILJ := 0..1,KM1 DO
186 BEGIN
187 KM := K + ILJ ;
188 LOOP I := K,K,L-K DO % LOOP NUMBER 1
189 BEGIN
190 IL := I .. ILJ ;
191 ODD1 := P(I+ILJ+PI);
192 P(P1+I+ILJ) := ODD1 + P(P1+J-IL) ;
193 P(P1+J-IL) := ODD1 - P(P1+J-IL) ;
194 END; % OF LOOP NUMBER 1
195
196 P(P1+L+ILJ) := P(P1+L+ILJ) + P(P1+L+ILJ) ;
197 IL := K - ILJ ;
198 P(P1+ILJ) := - P(P1+J-IL) - P(P1+J-IL) ;
199 ODD3 := P(P1+KM);
200 P(P1+KM) := -ODD3 - ODD3 ;
201 IL := 0 ;
202 LOOP I := KP,KP,L-KP DO % LOOP NUMBER 2
203 BEGIN
204 IL := IL + 2 ;
205 CJ := I + ILJ ;
206 ODD1 := P(P1+I+KM) - ODD3 ;
207 GL := IL.(16:42) ;
208 SMODE := MODE ; MODE := TRUE ;
209 RR := GRABONE(SGLJ,IL) ;
210 GL := (N2-IL).(16:42);
211 RRI := GRABONE(SGLJ,N2-IL);
212 MODE := SMODE ;
213 ODD2 := RR * P(P1+CJ) - RRI * ODD1 ;
214 ODD1 := RRI * P(P1+CJ) + RR * ODD1 ;
215 P(P1+CJ) := P(P1+J+KM-1) - P(P1+J-I-K+ILJ) ;
216 ODD3 := P(P1+I+KM) ;
217 P(P1+I+KM) := ODD2 - P(P1+J-I+ILJ);
218 P(P1+J-I+KM) := ODD2 + P(P1+J-I+ILJ);
219 P(P1+J-I+ILJ) := P(P1+CJ) - ODD1 ;
220 P(P1+CJ) := P(P1+CJ) + ODD1 ;
221 END ; % OF LOOP NUMBER 2
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00019900
0001 := P(P1+L+KM) ;
P(P1+L+KM) := P(P1+L+ILJ) ;
P(P1+L+ILJ) := 0001 + 0001 ;
LOOP I := ILJ,K,L-K+ILJ 00 % LOOP NUMBER 3
BEGIN
  0001 := P(P1+L) ;
  P(P1+L) := 0001 + P(P1+L+I) ;
  P(P1+L+I) := 0001 - P(P1+L+I) ;
  END ; % OF LOOP NUMBER 3
00022200
MCOE := TRUE ;
ENO ;
% SCRAMBLING
IL := 0 ;
KM := K+KM1 ;
LOOP I := KP,KP,J-KP DO
  BEGIN
    IL := IL + 1 ;
    GL := IL .(16:42) ;
    0003 := INDEX(GL) ;
    CJ := GRABONE(0003 ,IL) ;
    IF IL LSS CJ THEN
      BEGIN
        CJ := CJ * KP ;
        LOOP ILJ := 0,1,KM 00 % LOOP NUMBER 4
        BEGIN
          0001 := P( 1+ILJ) ;
          P(1+ILJ) := P (CJ+ILJ) ;
          P(CJ+ILJ) := 0001 ;
          END ; % OF LOOP NUMBER 4
        ENO ;
      ENO ;
    ENO ;
  IT := KP + KP ;
  IP := N3 ;
  LOOP I := 2,1,LLN -1 00
  BEGIN
    IP:=IP.(16:47) ;
    CL := 0 ;
    L2 := 1 ;
    CJ := IT.(16:47) ;
    IT := IT + IT ;
    LOOP IL := 0,CJ,CJ DO
      BEGIN
        LOOP KM := 1L,KP,IL+CJ-KP DO
          BEGIN
            GL := (N2-CL) .(16:42) ;
            CO := GRABONE(S(GL),N2-CL) * L2 ;
            GL := CL .(16:42) ;
            SI := GRABONE(S(GL),CL) ;
            LOOP L3 := 0,IT,J-IT 00
              BEGIN
                MODE := FMODE ;
                LOOP ILJ := 0,1,KM1 00 % LOOP NUMBER 5
                  BEGIN
                    L1 := L3 + KM + ILJ ;

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00021800 04:00:001682

00021900 04:00:001694

00022000 04:00:001716

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00022300 05:00:001775

00022400 05:00:001783

00022500 05:00:001801

00022600 05:00:001822

00022700 04:00:001830

00022800 04:00:001834

00018300

00022900 03:00:001842

00023000 03:00:001842

00023100 03:00:001847

00023200 03:00:001856

00023300 03:00:001902

00023400 04:00:001902

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00023700 04:00:001921

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00023900 04:00:001946

00024000 05:01:001952

00024100 05:01:001964

00024200 05:01:001988

00024300 06:01:001988

00024400 06:01:001997

00024500 06:01:002015

00024600 06:01:002024

00024700 05:01:002032

00023900 04:00:002034

00023300

00024900 03:00:002042

00025000 03:00:002051

00025100 03:00:002055

00025200 03:00:002082

00025300 04:00:002082

00025400 04:00:002088

00025500 04:00:002093

00025600 04:00:002098

00025700 04:00:002104

00025800 04:00:002113

00025900 04:00:002146

00026000 05:00:002146

00026100 05:00:002193

00026200 06:00:002193

00026300 06:00:002204

00026400 06:00:002232

00026500 06:00:002238

00026600 06:00:002249

00026700 06:00:002287

00026800 07:00:002287

00026900 07:00:002291

00027000 07:00:002315

00027100 08:00:002315

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272 L4 := LI + CJ + CJ ;
273 ODD1 := P(PI+L4) * CO - P(PI+L4+K) * SI ;
274 ODD2 := P(PI+L4+K) * CO + P(PI+L4) * SI ;
275 P(PI+L4) := P(PI+L1) - ODD1 ;
276 P(PI+L4+K) := P(PI+L1+K) - ODD2 ;
277 P(PI+L1) := P(PI+L1) + ODD1 ;
278 P(PI+L1+K) := P(PI+L1+K) + ODD2 ;
279 MODE := TRUE ;
280 END ; % OF LOOP NUMBER 5

281 END ;

282 CL := CL + IP ;
283 END ;

284 L2 := -I ;
285 IP := -IP ;
286 END ;

287 END ;

288 I := 0 ;
289 LOOP L1 := I, I+2 DO
290 BEGIN
291 I := I + K ;
292 GL := LI . [16:42] ;
293 PR := GRABONE(ISS(GL), L1) ;
294 MODE := FMODE ;
295 LOOP ILJ := 0, 1, KM1 DO % LOOP NUMBER 6
296 BEGIN
297 ODD1 := P(PI+I+ILJ) - P(PI+J-I+ILJ) ;
298 ODD2 := ( P(PI+I+ILJ) + P(PI+J-I+ILJ) ) * RR ;
299 P(PI+I+ILJ) := ODD2 + ODD1 ;
300 P(PI+J-I+ILJ) := ODD2 - ODD1 ;
301 MODE := TRUE ;
302 END ; % OF LOOP NUMBER 6

303 END ;

304 %
305 END OF FOURIER ANALYSIS AND SYNTHESIS
306 IF (I=2 THEN GO TO FINISH ;
307 %
308 % SECTION 3 %
309 % SKEW POTENTIAL
310 %
311 I := 0 ;
312 LOOP CJ := KP, K, KC DO % LOOP NUMBER 7
313 BEGIN
314 I := I + I ;
315 LOOP IL := CJ, 1, CJ+KM1 DO
316 BEGIN
317 P(IL) := RTR(I, P(IL)) ;
318 END ;
319 END ; % OF LOOP NUMBER 7

320 END ; % OF LOOP NUMBER 7

321 CRED STARTS HERE
322 MODE := (PEN LSS CN-I) ;
323 PI := PEN * K ;
324 SINCE P(0) TO P(KM1) IS ALREADY GARBAGE WE WONT PROTECT IT
325 CNIK := CN.[16:42] - 1 ; IF CNIK LSS 0 THEN CNIK:=0 ;
326 CJ := CM.[158:6] ;

```

```

00027200 09:00:002324
00027300 08:00:002336
00027400 08:00:002365
00027500 08:00:002393
00027600 08:00:002410
00027700 08:00:002433
00027800 08:00:002450
00027900 08:00:002473
00028000 08:00:002477
00027000
00028100 07:00:002485
00026700 06:00:002493
00028300 06:00:002502
00026100
00028400 05:00:002510
00028500 05:00:002516
00028600 05:00:002522
00025900
00028700 04:00:002530
00025200
00028800 03:00:002538
00028900 03:00:002543
00029000 03:00:002567
00029100 04:00:002567
00029200 04:00:002576
00029300 04:00:002581
00029400 04:00:002591
00029500 04:00:002595
00029600 04:00:002619
00029700 05:00:002619
00029800 05:00:002645
00029900 05:00:002674
00030000 05:00:002686
00030100 05:00:002701
00030200 05:00:002705
00029600
00030300 04:00:002713
00029000
00030400 03:00:002721
00030500 03:00:002721
00030600 03:00:002742
00030700 03:00:002742
00030800 03:00:002742
00030900 03:00:002747
00031000 03:00:002788
00031100 04:00:002788
00031200 04:00:002795
00031300 04:00:002830
00031400 05:00:002830
00031500 05:00:002846
00031300
00031600 04:00:002854
00031000
00031700 03:00:002862
00031800 03:00:002862
00031900 03:00:002876
00032000 03:00:002884
00032100 03:00:002884
00032200 03:00:002917

```



```

323 IF CJ=0 THEN CJ := 64 ;
324 LOOP I := 0,1,CNIK DO
325 BEGIN
326   IL := 4 ;
327   L4 := 0 ;
328   E := 1.0 / A[I] ;
329   Q := P[PI] * E ;
330   CL := 64 ;
331   LOOP L1 := 0,1,KM1 DO
332   BEGIN
333     IF L1=KM1 THEN CL := CJ ;
334     LOOP L2 := IL,1,CL DO % LOOP NUMBER 8
335     BEGIN
336       MODE := REVR(1,MODE) ;
337       A[I] := RTR(1,,A[I]) ;
338       E := 1.0 / ( A[I] - PTR(1,,E) ) ;
339       PI := RTR(1,,PI) ;
340       Q := (P[PI+L1+L4] - RTR(1,,Q)) * E ;
341       L4 := 0 ;
342     END % OF LOOP NUMBER 8
343
344     L4 := -1 ;
345     IL := 1 ;
346   END ;
347
348   CL := 64 ;
349   E := P[PI+KM1] - A[I] * Q ;
350   P[PI+KM1] := Q ;
351   IL := 4 ;
352   L4 := 0 ;
353   LOOP L1 := 0,1,KM1 DO
354   BEGIN
355     IF L1=KM1 THEN CL := CJ ;
356     L3 := KM1 - L1 ;
357     LOOP L2 := L1,1,CL DO % LOOP NUMBER 9
358     BEGIN
359       MODE := REVL(1,MODE) ;
360       PI := RTR(1,,PI) ;
361       A[I] := RTR(1,,A[I]) ;
362       Q := RTR(1,,Q) ;
363       E := RTR(1,,E) ;
364       T := P[PI+L3+L4] - Q - A[I] * E ;
365       Q := E ;
366       E := T ;
367       P[PI+L3+L4] := Q ;
368       L4 := 0 ;
369     END % OF LOOP NUMBER 9
370
371     IL := 1 ; L4 := 1 ;
372   END ;
373
374   PI := PI + 64 * K ;
375   END ;
376
377   MODE := TRUE ; % UNSKEW POTENTIAL
378   I := 0 ;
379   LOOP CJ := KP,K,KC DO
380   BEGIN
381     I := I + 1 ;
382     LOOP IL := CJ,1,CJ+KM1 DO % LOOP NUMBER 10
383
00032300 03:00:002923
00032400 03:00:002945
00022503 03:00:002970
00032600 04:00:002970
00032700 04:00:002975
00032803 04:00:002980
00032900 04:00:002990
00033000 04:00:002995
00033100 04:00:003000
00033200 04:00:003024
00033300 05:00:003024
00033400 05:00:003047
00033500 05:00:003079
00033600 06:00:003079
00033700 06:00:003083
00033800 06:00:003097
00033900 06:00:003119
00034000 06:00:003127
00034100 06:00:003149
00034200 06:00:003154
00033500
00034300 05:00:003162
00034400 05:00:003168
00034500 05:00:003173
00033200
00034600 04:00:003181
00034700 04:00:003186
00034800 04:00:003202
00034900 04:00:003210
00035000 04:00:003215
00035100 04:00:003220
00035200 04:00:003244
00035300 05:00:003244
00035400 05:00:003267
00035500 05:00:003275
00035600 05:00:003307
00035700 06:00:003307
00035800 06:00:003311
00035900 06:00:003319
00036003 06:00:003333
00036100 06:00:003341
00036200 06:00:003349
00036300 06:00:003370
00036400 06:00:003372
00036500 06:00:003374
00036600 06:00:003385
00036700 06:00:003390
00035600
00036800 05:00:003398
00036900 05:00:003408
00035200
00037000 04:00:003416
00037100 04:00:003425
00032500
00037200 03:00:003433
00037300 03:00:003437
00037400 03:00:003442
00037500 03:00:003483
00037600 04:00:003483
00037700 04:00:003490

```

```

378 BEGIN
379 P(I) := RTL(I,P(I)) ;
380 END: % OF LOOP NUMBER 10

381 END;

382 GO TO FOURIER ;
383 ***** SECTION 4 *****
384 FINISH: LOOP 1 := 0,1,KM DO
385 BEGIN
386 C(I+K) := STORE(I) ;
387 C(I+K) := STORE(I+K) ;
388 P(I) := STORE(I+K) ;
389 END ;

390 MODE := SAVEMODE ;
391 END ;

392 *****
393 ***** A TEST PROGRAM *****
394 CINT CN,CM,PP,I,J,CK,CS,OFF ;
395 CREAL HX,HY ;
396 PREAL VECTOR INDEX(I) ;
397 PREAL VECTOR AX,SA,ISX(I) ; PREAL VECTOR P, C(30) ;
398 PREAL X ;
399 SUBROUTINE DATA;
400 % THIS IS A DUMMY SUBROUTINE USED TO INITIALIZE P,
401 % THE POTENTIAL.
402 CODE
403 USE P ;
404 BEGIN
405 HERE:
406 SET;
407 ORG P;
408 DATA
409 0.50337,0.93489,0.60033,0.44270,0.10798,0.92968,0.92879,0.81380,(0.0156,
410 0.98078,0.52103,0.21337,0.52409,0.60700,0.30343,0.14217,0.33953,(0.0156,
411 0.79932,0.11226,0.55894,0.10389,0.40150,0.79314,0.31545,0.69259,(0.0156,
412 0.34400,0.37663,0.58489,0.45073,0.14357,0.52035,0.12551,0.19098,(0.0156,
413 0.63619,0.06807,0.78106,0.75710,0.66561,0.86559,0.40329,0.02445,(0.0156,
414 0.15432,0.90765,0.08608,0.18121,0.37631,0.54454,0.24602,0.73611,(0.0156,
415 0.95768,0.5112,0.80428,0.75255,0.07879,0.16337,0.07416,0.56654,(0.0156,
416 0.08343,0.76303,0.43107,0.60956,0.50404,0.02133,0.075174,0.97407,(0.0156;
417 ORG HERE;
418 END CODE;
419 *****
420 CN := 9;
421 CM := 8 ;
422 PP:= 3 ;
423 CK := CN - 1 ;
424 HX := CK * CK ;
425 HY := (CM-1)*(CM-1) ;
426 CS := CK ;
427 SETPT(AX,SA,ISX,PP,INDEX,CK,HX,HY) ;
428 LOOP I := 1,1,CS-1 DO BEGIN
429 C(I) := HX*(P(I-1)+P(I+1))-2*P(I)+HY*(RTL(I)+P(I))
430 +RTL(1,P(I))-2*P(I)) ;
431 END ;

432 LOOP I:=0,1,CS DO SIMWRITE(LINE,"BP(Y)",P(I)) ;

```

```

00037800 04:00:003525
00037900 05:00:003525
00038000 05:00:003544
00037800
00037500
00038200 03:00:003560
00038300 03:00:003561
00038400 03:00:003561
00038500 03:00:003586
00038600 04:00:003586
00038700 04:00:003500
00038800 04:00:003519
00038900 04:00:003632
00038500
00039000 03:00:003640
00039100 03:00:003644
00009600
00039200 01:00:003662
00039300 01:00:003662
00039400 01:00:003662
00039500 01:00:003662
00039600 01:00:003662
00039700 01:00:003662
00039800 01:00:003662
00039900 01:00:003662
00040000 02:00:003662
00040100 02:00:003662
00040200 02:00:003662
00040300 02:00:003662
00040400 02:00:003663
00040500 02:00:003663
00040600 02:00:003663
00040700 02:00:003663
00040800 02:00:003663
00040900 02:00:003663
00041000 02:00:003663
00041100 02:00:003663
00041200 02:00:003663
00041300 02:00:003663
00041400 02:00:003663
00041500 02:00:003663
00041600 02:00:003663
00041700 02:00:003663
00041800 02:00:003663
00041900 01:00:003675
00042000 01:00:003675
00042100 01:00:003680
00042200 01:00:003685
00042300 01:00:003690
00042400 01:00:003697
00042500 01:00:003714
00042600 01:00:003735
00042700 01:00:003739
00042800 01:00:003770
00042900 02:00:003797
00043000 02:00:003849
00043100 02:00:003873
00042800
00043200 01:00:003881

```

```

433 POISSONDIRECT(P,C,AX,SX,CN,CM,HX,HY      ,ISX,PP,INDEX);
434 LOOP I := 0,1,CS DO BEGIN
435     SIMPITE(LINE,"XP(I)",I,PI));
436 END ;
437 END.

```

```

00043300 01:00:003933
00043400 01:00:003968
00043500 02:00:003992
00043600 02:00:004022
00043400
00043700 01:00:004030
00000100

```

NUMBER OF ERRORS DETECTED=0000. NUMBER OF WARNINGS NOTED=0000.  
 SOURCE PROGRAM SIZE= 00437 CARD IMAGES, 004209 SYNTACTIC ITEMS.  
 SOURCE FILE: CARD=ERICKSEN/PRINT/HOCKNEYG.  
 MACRO LIBRARY: ILLIAC/GLYPNIR3/MACROS.  
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